

chain nodes :

6 9 13 14 15 16 19 22 23 24 26

ring nodes :

1 2 3 4 5 27 28 29 30 31 32 33 34 35 36

chain bonds :

2-26 5-6 13-14 15-16 22-23 22-24

ring bonds :

1-2 1-5 2-3 3-4 4-5 27-28 27-32 28-29 29-30 30-31 30-33 31-32
31-36 33-34 34-35 35-36

exact/norm bonds :

1-2 1-5 2-3 2-26 3-4 4-5 5-6 13-14 15-16 22-23 22-24 27-28
27-32 28-29 29-30 30-31 30-33 31-32 31-36 33-34 34-35 35-36

isolated ring systems :

containing 1 :

G1:CH2,O,[*1]

G2:[*2],[*3],[*4],[*5]

G3:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 9:CLASS 13:CLASS
14:CLASS 15:CLASS 16:CLASS 19:CLASS 22:CLASS 23:CLASS 24:CLASS
26:CLASS 27:CLASS 28:CLASS 29:CLASS 30:Atom 31:Atom 32:Atom 33:CLASS
34:CLASS 35:CLASS 36:CLASS

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=> d his

(FILE 'HOME' ENTERED AT 16:43:46 ON 25 JAN 2007)

FILE 'REGISTRY' ENTERED AT 16:43:54 ON 25 JAN 2007

L1 STRUCTURE UPLOADED

L2 0 S L1

L3 37 S L1 FULL

FILE 'HCAPLUS' ENTERED AT 16:53:32 ON 25 JAN 2007

L4 1 S L3

FILE 'CAOLD' ENTERED AT 16:53:59 ON 25 JAN 2007

L5 0 S L3

FILE 'REGISTRY' ENTERED AT 16:54:06 ON 25 JAN 2007

L6 STRUCTURE UPLOADED

L7 0 S L6

L8 STRUCTURE UPLOADED

L9 0 S L8

L10 STRUCTURE UPLOADED

L11 0 S L10

L12 STRUCTURE UPLOADED

L13 0 S L12

L14 35 S L12 FULL

FILE 'HCAPLUS' ENTERED AT 17:09:54 ON 25 JAN 2007

L15 1 S L14

FILE 'REGISTRY' ENTERED AT 17:10:06 ON 25 JAN 2007

L16 STRUCTURE UPLOADED

L17 0 S L16

L18 STRUCTURE UPLOADED

L19 0 S L18

L20 STRUCTURE UPLOADED

L21 0 S L20

L22 0 S L20 FULL

L23 35 S L20 FULL

L24 0 S L23 NOT L14

FILE 'REGISTRY' ENTERED AT 17:15:33 ON 25 JAN 2007

L25 STRUCTURE UPLOADED

L26 0 S L25

L27 STRUCTURE UPLOADED

L28 0 S L27

L29 35 S L27 FULL

L30 STRUCTURE UPLOADED

L31 0 S L30

L32 35 S L30 FULL

L33 STRUCTURE UPLOADED

L34 STRUCTURE UPLOADED

L35 0 S L34

L36 6 S L34 FULL

FILE 'HCAPLUS' ENTERED AT 17:24:05 ON 25 JAN 2007

L37 5 S L36

FILE 'CAOLD' ENTERED AT 17:24:31 ON 25 JAN 2007

Updated Search

4 1)

10508761

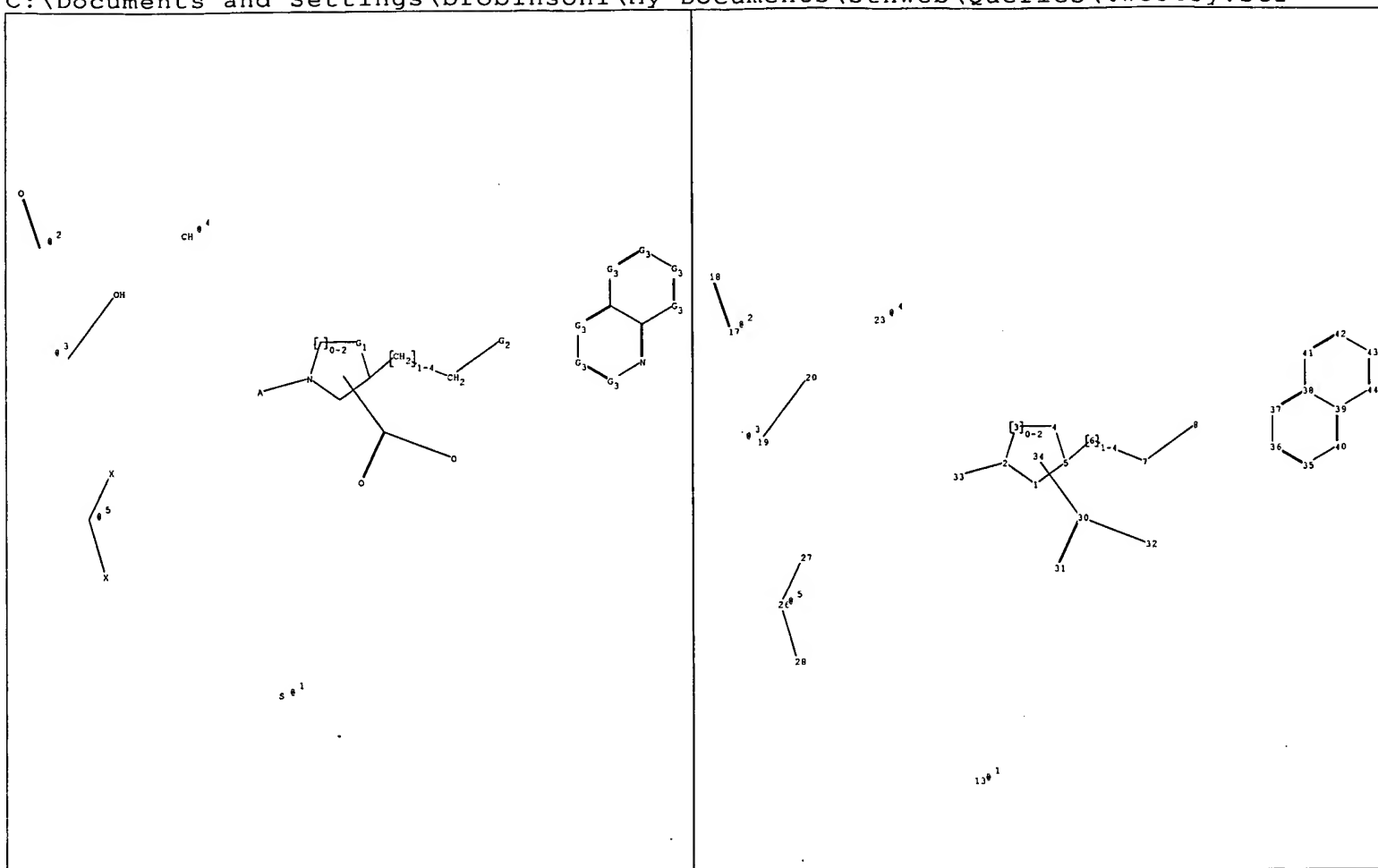
=> s 136
L38 0 L36

=> s 123
L39 , 0 L23

=>

Updated Search





chain nodes :

6 7 8 13 17 18 19 20 23 26 27 28 30 31 32 33

ring nodes :

1 2 3 4 5 35 36 37 38 39 40 41 42 43 44

chain bonds :

2-33 5-6 6-7 7-8 17-18 19-20 26-27 26-28 30-32 30-31

ring bonds :

1-2 1-5 2-3 3-4 4-5 35-36 35-40 36-37 37-38 38-39 38-41 39-40
39-44 41-42 42-43 43-44

exact/norm bonds :

1-2 1-5 2-3 2-33 3-4 4-5 5-6 6-7 7-8 17-18 19-20 26-27 26-28
30-32 30-31 35-36 35-40 36-37 37-38 38-39 38-41 39-40 39-44 41-42
42-43 43-44

isolated ring systems :

containing 1 :

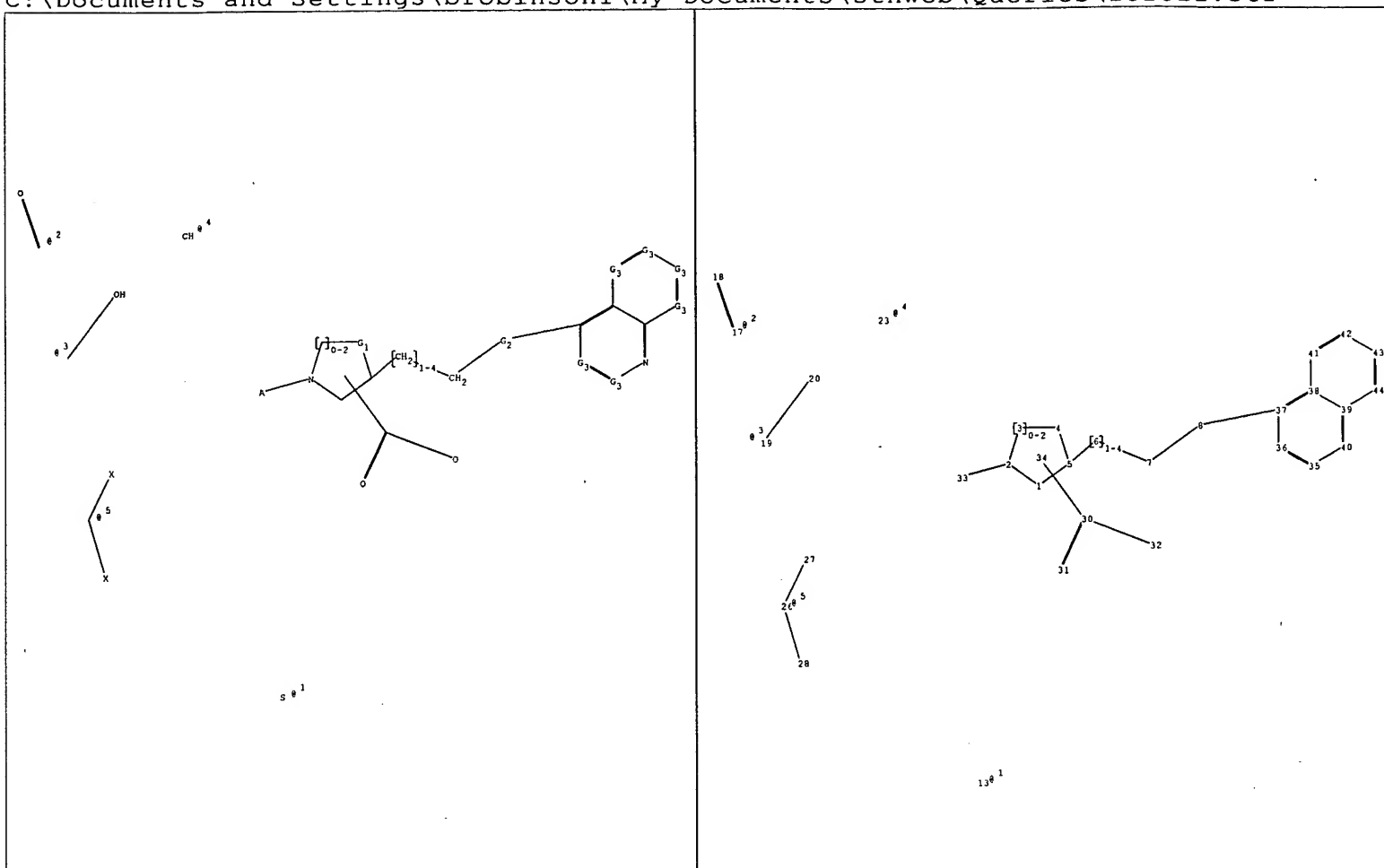
G1:CH2,O,[*1]

G2:[*2],[*3],[*4],[*5]

G3:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:CLASS 8:CLASS 13:CLASS
17:CLASS 18:CLASS 19:CLASS 20:CLASS 23:CLASS 26:CLASS 27:CLASS
28:CLASS 30:CLASS 31:CLASS 32:CLASS 33:CLASS 34:CLASS 35:CLASS
36:CLASS 37:CLASS 38:Atom 39:Atom 40:Atom 41:CLASS 42:CLASS 43:CLASS
44:CLASS



chain nodes :

6 7 8 13 17 18 19 20 23 26 27 28 30 31 32 33

ring nodes :

1 2 3 4 5 35 36 37 38 39 40 41 42 43 44

chain bonds :

2-33 5-6 6-7 7-8 8-37 17-18 19-20 26-27 26-28 30-32 30-31

ring bonds :

1-2 1-5 2-3 3-4 4-5 35-36 35-40 36-37 37-38 38-39 38-41 39-40
39-44 41-42 42-43 43-44

exact/norm bonds :

1-2 1-5 2-3 2-33 3-4 4-5 5-6 6-7 7-8 8-37 17-18 19-20 26-27
26-28 30-32 30-31 35-36 35-40 36-37 37-38 38-39 38-41 39-40 39-44
41-42 42-43 43-44

isolated ring systems :

containing 1 :

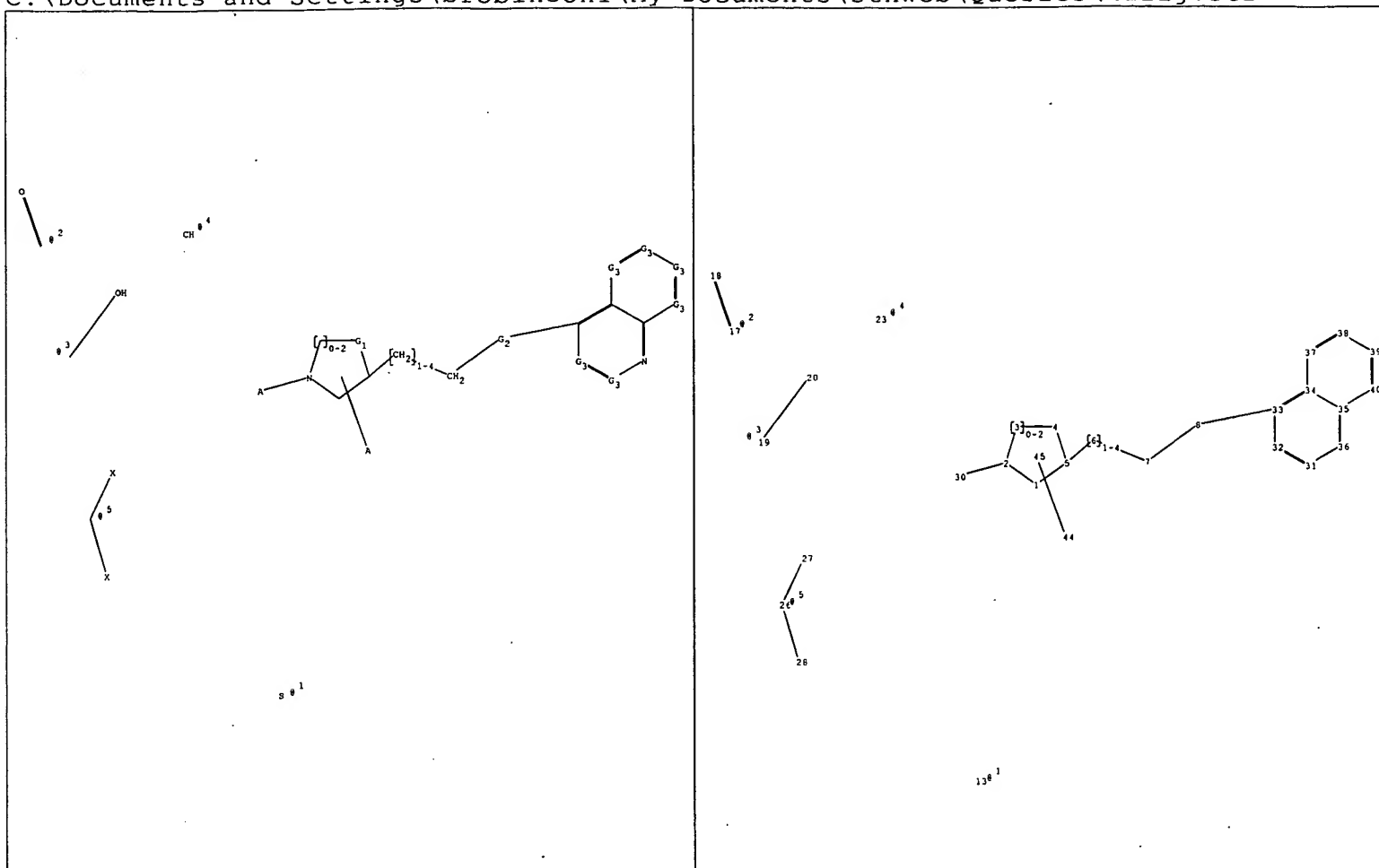
G1:CH2,O,[*1]

G2:[*2],[*3],[*4],[*5]

G3:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:CLASS 8:CLASS 13:CLASS
17:CLASS 18:CLASS 19:CLASS 20:CLASS 23:CLASS 26:CLASS 27:CLASS
28:CLASS 30:CLASS 31:CLASS 32:CLASS 33:CLASS 34:CLASS 35:CLASS
36:CLASS 37:CLASS 38:Atom 39:Atom 40:Atom 41:CLASS 42:CLASS 43:CLASS
44:CLASS



chain nodes :

6 7 8 13 17 18 19 20 23 26 27 28 30 44

ring nodes :

1 2 3 4 5 31 32 33 34 35 36 37 38 39 40

chain bonds :

2-30 5-6 6-7 7-8 8-33 17-18 19-20 26-27 26-28

ring bonds :

1-2 1-5 2-3 3-4 4-5 31-32 31-36 32-33 33-34 34-35 34-37 35-36
35-40 37-38 38-39 39-40

exact/norm bonds :

1-2 1-5 2-3 2-30 3-4 4-5 5-6 6-7 7-8 8-33 17-18 19-20 26-27
26-28 31-32 31-36 32-33 33-34 34-35 34-37 35-36 35-40 37-38 38-39
39-40

isolated ring systems :

containing 1 :

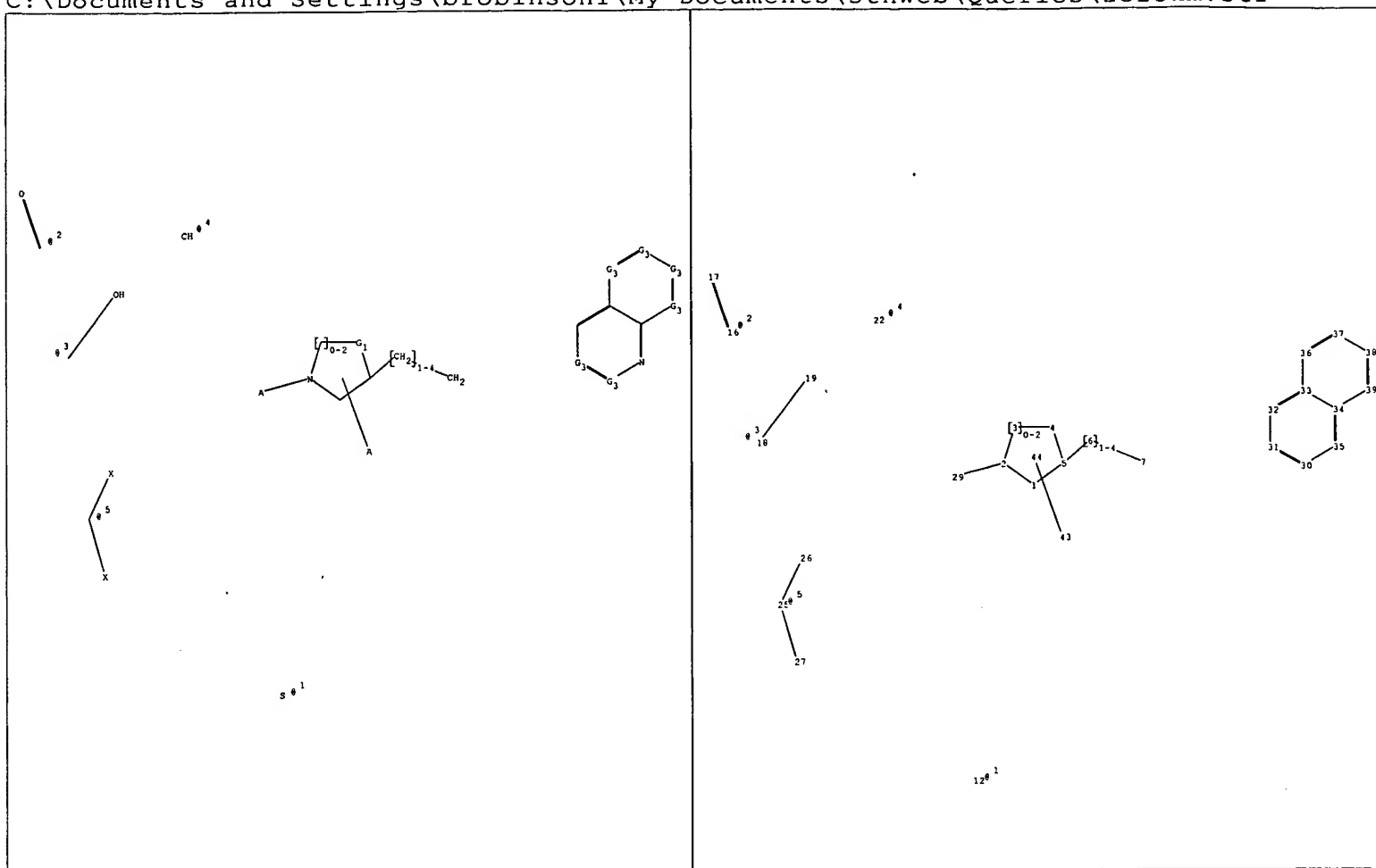
G1:CH2,O,[*1]

G2:[*2],[*3],[*4],[*5]

G3:C,N

Match level :

1:Atom	2:Atom	3:Atom	4:Atom	5:Atom	6:CLASS	7:CLASS	8:CLASS	13:CLASS
17:CLASS	18:CLASS	19:CLASS	20:CLASS	23:CLASS	26:CLASS	27:CLASS		
28:CLASS	30:CLASS	31:CLASS	32:CLASS	33:CLASS	34:Atom	35:Atom	36:Atom	
37:CLASS	38:CLASS	39:CLASS	40:CLASS	44:CLASS	45:CLASS			



chain nodes :

6 7 12 16 17 18 19 22 25 26 27 29 43

ring nodes :

1 2 3 4 5 30 31 32 33 34 35 36 37 38 39

chain bonds :

2-29 5-6 6-7 16-17 18-19 25-26 25-27

ring bonds :

1-2 1-5 2-3 3-4 4-5 30-31 30-35 31-32 32-33 33-34 33-36 34-35
34-39 36-37 37-38 38-39

exact/norm bonds :

1-2 1-5 2-3 2-29 3-4 4-5 5-6 6-7 16-17 18-19 25-26 25-27 30-31
30-35 31-32 32-33 33-34 33-36 34-35 34-39 36-37 37-38 38-39

isolated ring systems :

containing 1 :

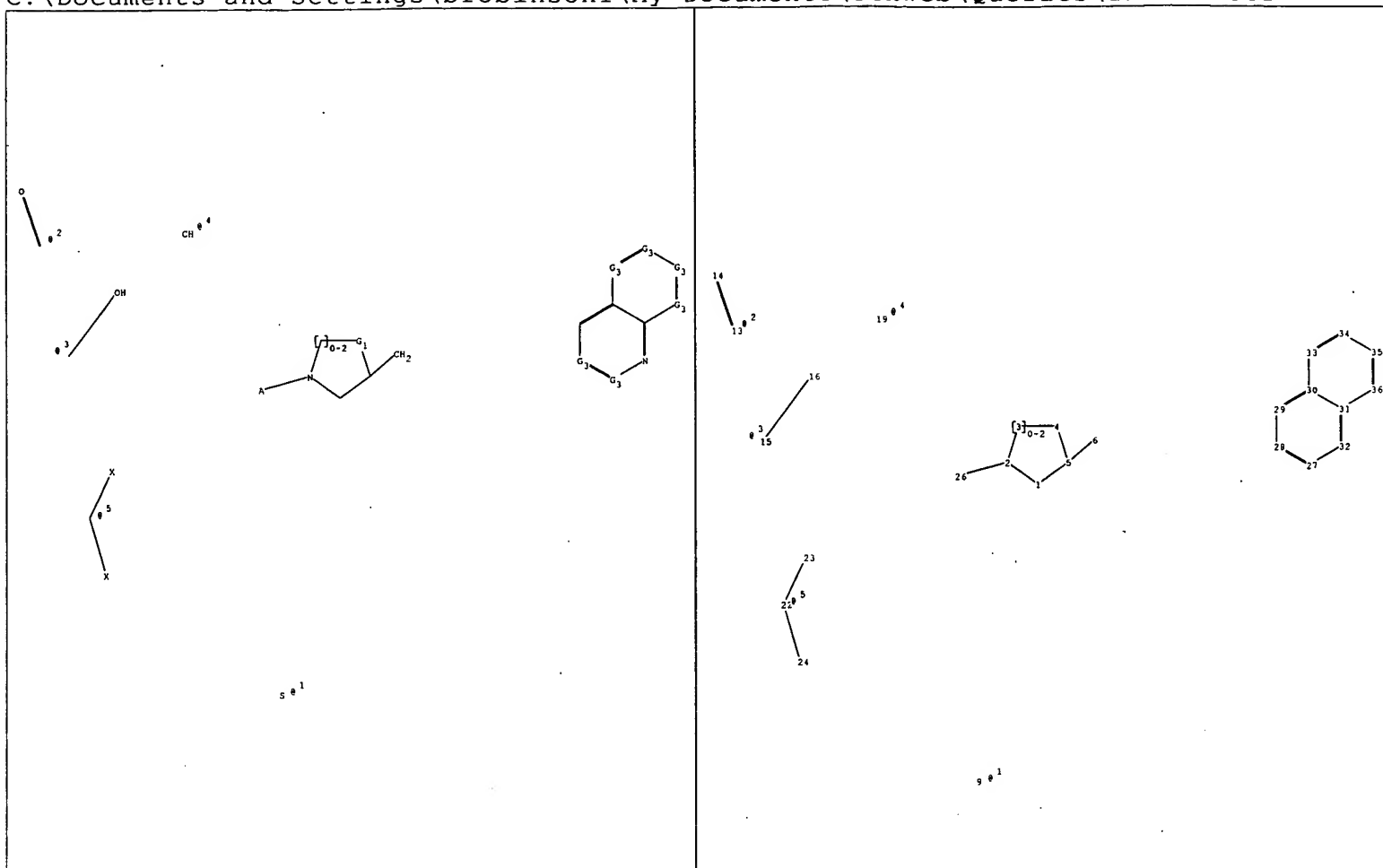
G1:CH2,O,[*1]

G2:[*2],[*3],[*4],[*5]

G3:C,N

Match level :

1:Atom	2:Atom	3:Atom	4:Atom	5:Atom	6:CLASS	7:CLASS	12:CLASS
16:CLASS	17:CLASS	18:CLASS	19:CLASS	22:CLASS	25:CLASS	26:CLASS	
27:CLASS	29:CLASS	30:CLASS	31:CLASS	32:CLASS	33:Atom	34:Atom	35:Atom
36:CLASS	37:CLASS	38:CLASS	39:CLASS	43:CLASS	44:CLASS		



chain nodes :

6 9 13 14 15 16 19 22 23 24 26

ring nodes :

1 2 3 4 5 27 28 29 30 31 32 33 34 35 36

chain bonds :

2-26 5-6 13-14 15-16 22-23 22-24

ring bonds :

1-2 1-5 2-3 3-4 4-5 27-28 27-32 28-29 29-30 30-31 30-33 31-32
31-36 33-34 34-35 35-36

exact/norm bonds :

1-2 1-5 2-3 2-26 3-4 4-5 5-6 13-14 15-16 22-23 22-24 27-28
27-32 28-29 29-30 30-31 30-33 31-32 31-36 33-34 34-35 35-36

isolated ring systems :

containing 1 :

G1:CH2,O,[*1]

G2:[*2],[*3],[*4],[*5]

G3:C,N

Match level :

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14:CLASS 15:CLASS 16:CLASS 19:CLASS 22:CLASS 23:CLASS 24:CLASS
26:CLASS 27:CLASS 28:CLASS 29:CLASS 30:Atom 31:Atom 32:Atom 33:CLASS
34:CLASS 35:CLASS 36:CLASS

10508761

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal612bxr

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 "Ask CAS" for self-help around the clock
NEWS 3 OCT 23 The Derwent World Patents Index suite of databases on STN
has been enhanced and reloaded
NEWS 4 OCT 30 CHEMLIST enhanced with new search and display field
NEWS 5 NOV 03 JAPIO enhanced with IPC 8 features and functionality
NEWS 6 NOV 10 CA/CAPLUS F-Term thesaurus enhanced
NEWS 7 NOV 10 STN Express with Discover! free maintenance release Version
8.01c now available
NEWS 8 NOV 20 CAS Registry Number crossover limit increased to 300,000 in
additional databases
NEWS 9 NOV 20 CA/CAPLUS to MARPAT accession number crossover limit increased
to 50,000
NEWS 10 DEC 01 CAS REGISTRY updated with new ambiguity codes
NEWS 11 DEC 11 CAS REGISTRY chemical nomenclature enhanced
NEWS 12 DEC 14 WPIDS/WPINDEX/WPIX manual codes updated
NEWS 13 DEC 14 GBFULL and FRFULL enhanced with IPC 8 features and
functionality
NEWS 14 DEC 18 CA/CAPLUS pre-1967 chemical substance index entries enhanced
with preparation role
NEWS 15 DEC 18 CA/CAPLUS patent kind codes updated
NEWS 16 DEC 18 MARPAT to CA/CAPLUS accession number crossover limit increased
to 50,000
NEWS 17 DEC 18 MEDLINE updated in preparation for 2007 reload
NEWS 18 DEC 27 CA/CAPLUS enhanced with more pre-1907 records
NEWS 19 JAN 08 CHEMLIST enhanced with New Zealand Inventory of Chemicals
NEWS 20 JAN 16 CA/CAPLUS Company Name Thesaurus enhanced and reloaded
NEWS 21 JAN 16 IPC version 2007.01 thesaurus available on STN
NEWS 22 JAN 16 WPIDS/WPINDEX/WPIX enhanced with IPC 8 reclassification data
NEWS 23 JAN 22 CA/CAPLUS updated with revised CAS roles
NEWS 24 JAN 22 CA/CAPLUS enhanced with patent applications from India

NEWS EXPRESS NOVEMBER 10 CURRENT WINDOWS VERSION IS V8.01c, CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
AND CURRENT DISCOVER FILE IS DATED 25 SEPTEMBER 2006.

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=>

=> file reg

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.21	0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 16:43:54 ON 25 JAN 2007

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DICTIONARY FILE UPDATES: 24 JAN 2007 HIGHEST RN 918400-64-3

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=>

Uploading C:\Documents and Settings\brobinson1\My Documents\stnweb\Queries\987tys.str

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 16:53:24 FILE 'REGISTRY'

Updated Search

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SAMPLE SCREEN SEARCH COMPLETED - 467 TO ITERATE

100.0% PROCESSED 467 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 8044 TO 10636
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 full

THE ESTIMATED SEARCH COST FOR FILE 'REGISTRY' IS 171.65 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y
FULL SEARCH INITIATED 16:53:29 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 9164 TO ITERATE

100.0% PROCESSED 9164 ITERATIONS 37 ANSWERS
SEARCH TIME: 00.00.01

L3 37 SEA SSS FUL L1

=> file hcaplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	178.85	179.06

FILE 'HCAPLUS' ENTERED AT 16:53:32 ON 25 JAN 2007.
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FILE COVERS 1907 - 25 Jan 2007 VOL 146 ISS 5
FILE LAST UPDATED: 24 Jan 2007 (20070124/ED)

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l3

L4 1 L3

=> d l4, ibib abs hitstr, 1

L4 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2007 ACS on STN
ACCESSION NUMBER: 2004:800851 HCAPLUS

Updated Search

10508761

DOCUMENT NUMBER: 141:314170
 TITLE: 4-Substituted quinoline derivatives, the preparation thereof and compositions containing same, useful as antimicrobials
 INVENTOR(S): Bigot, Antony; El Ahmad, Youssef; Malleron, Jean Luc; Martin, Jean Paul; Mignani, Serge; Pantel, Guy; Ronan, Baptiste; Tabart, Michel
 PATENT ASSIGNEE(S): Aventis Pharma SA, Fr.
 SOURCE: Fr. Demande, 67 pp.
 CODEN: FRXXBL
 DOCUMENT TYPE: Patent
 LANGUAGE: French
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
FR 2852954	A1	20041001	FR 2003-3812	20030328
FR 2852954	B1	20060714		
US 2004224946	A1	20041111	US 2004-810711	20040326
AU 2004226207	A1	20041014	AU 2004-226207	20040329
CA 2520764	A1	20041014	CA 2004-2520764	20040329
WO 2004087647	A2	20041014	WO 2004-FR783	20040329
WO 2004087647	A3	20050127		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
EP 1611127	A2	20060104	EP 2004-742385	20040329
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK				
CN 1795191	A	20060628	CN 2004-80014510	20040329
JP 2006522779	T	20061005	JP 2006-505763	20040329
PRIORITY APPLN. INFO.:			FR 2003-3812	A 20030328
			US 2003-487084P	P 20030714
			WO 2004-FR783	W 20040329

OTHER SOURCE(S): MARPAT 141:314170
 GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB Quinoline-4-substituted derivs. I are disclosed [wherein X, Y, Z, U, T = C-R1' to CR5' resp., or one or more is a N atom; R1, R1', R2', R3', R4', R5' = independently H, halo, cyclo/alkyl, Ph, phenylthio, mono or bicyclic hetero(aryl)thio, OH and derivs., SH and derivs., NH2 and derivs., acyl, OCF3, OCHF2, CN, CO2H and derivs., NO2, etc.; D = CHR, CO, CROH, CRF, CF2; R = H, alkyl; A = (CH2)m; m = 1-3; B = (CH2)n; n = 0-2; E = CH2, and when Z = O, S, SO, SO2, then n = 2; R2 = CO2R, CH2CH2CO2R, CH2OH, CH2CH2OH, where R is defined as above; R3 = Ph, mono or bicyclic heteroaryl,

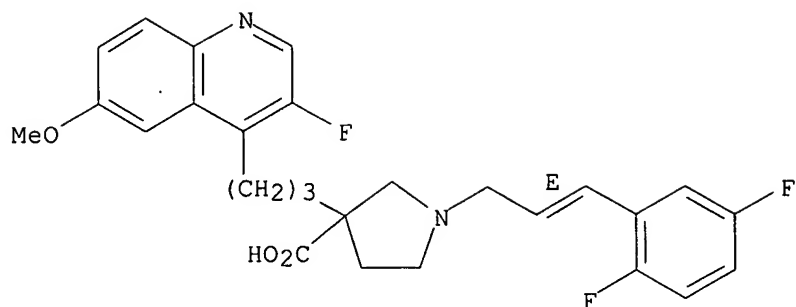
Updated Search

alkylene-R3'', etc.; R3'' = H, halo, OH and derivs., alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, cycloalkyl, acyl, Ph, OPh, heteroaryloxy, mono and bicyclic heteroaryl, NH2 and derivs., CONH2 and derivs., etc.; their enantiomers or diastereoisomers or their mixts., and/or their syn or anti forms or their mixts.; and their salts]. The novel derivs. are particularly interesting as antimicrobial agents. For example, II was prepared by amination of 2-[(E)-3-chloro-1-propenyl]-1,4-difluorobenzene (preparation given) with amine salt III•2HCl, followed by acidic hydrolysis. Compds. I were active against exptl. infections of mice by Staphylococcus aureus IP 8203 at 5-50 mg/kg s.c. or orally. None of the compds. showed toxicity in mice at 50 mg/kg s.c. (2 administrations).

IT 767355-23-7P, 1-[(E)-3-(2,5-Difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 767355-29-3P, (+)-1-[(E)-3-(2,5-Difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 767355-31-7P, (-)-1-[(E)-3-(2,5-Difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 767355-33-9P, (-)-2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[2-[(2,5-difluorophenyl)sulfanyl]ethyl]morpholine-2-carboxylic acid
 767355-35-1P, (+)-2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[2-[(2,5-difluorophenyl)sulfanyl]ethyl]morpholine-2-carboxylic acid
 767355-42-0P, (-)-2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[2-(2,5-difluorophenyl)allyl]morpholine-2-carboxylic acid
 767355-44-2P, (+)-2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[2-(2,5-difluorophenyl)allyl]morpholine-2-carboxylic acid
 767355-47-5P, 1-[(E)-3-(2,5-Difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]azetidine-3-carboxylic acid
 767355-52-2P, 3-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-1-[2-[(thiophen-2-yl)sulfanyl]ethyl]azetidine-3-carboxylic acid sodium salt
 767355-56-6P, 1-[2-(2,5-Difluorophenyl)sulfanyl]ethyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 767355-57-7P, 1-[2-(2,5-Difluorophenyloxy)ethyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 767355-58-8P, 1-[2-(Thiophen-2-ylsulfanyl)ethyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 767355-60-2P, 1-[(E)-3-(2,5-Difluorophenyl)-2-propenyl]-3-[3-(3-chloro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 767355-61-3P, 1-[2-(2,5-Difluorophenyl)sulfanyl]ethyl]-3-[3-(3-chloro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 767355-62-4P, 1-[2-(2,5-Difluorophenyloxy)ethyl]-3-[3-(3-chloro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 767355-63-5P, 1-[2-(Thiophen-2-ylsulfanyl)ethyl]-3-[3-(3-chloro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylic acid
 RL: ADV (Adverse effect, including toxicity); PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
 (bactericide; preparation of 4-substituted quinolines as antimicrobials)
 RN 767355-23-7 HCAPLUS
 CN 3-Pyrrolidinecarboxylic acid, 1-[(2E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxy-4-quinoliny)propyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

10508761

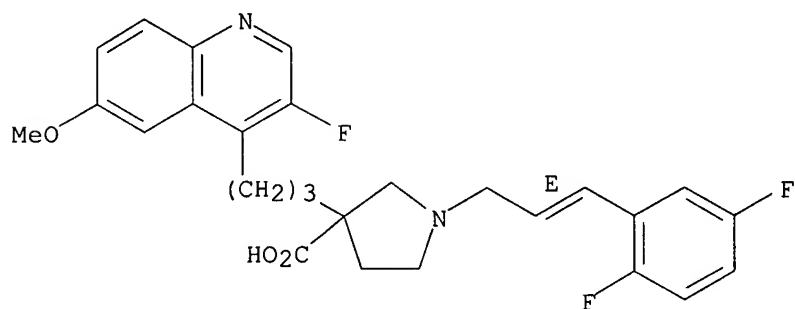


RN 767355-29-3 HCAPLUS

CN 3-Pyrrolidinecarboxylic acid, 1-[(2E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, (+)- (9CI) (CA INDEX NAME)

Rotation (+).

Double bond geometry as shown.

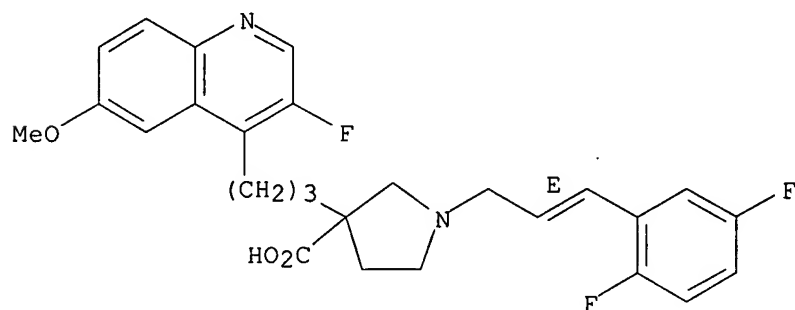


RN 767355-31-7 HCAPLUS

CN 3-Pyrrolidinecarboxylic acid, 1-[(2E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, (-)- (9CI) (CA INDEX NAME)

Rotation (-).

Double bond geometry as shown.



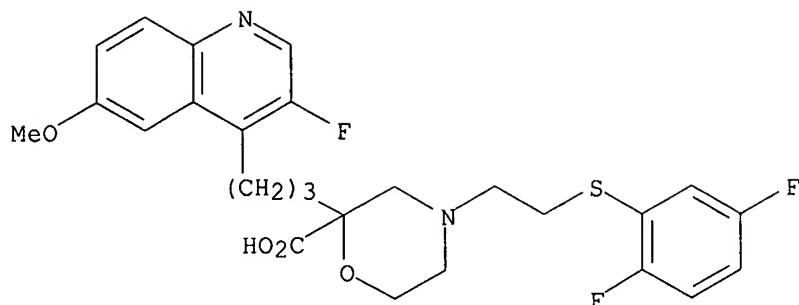
RN 767355-33-9 HCAPLUS

CN 2-Morpholinecarboxylic acid, 4-[2-[(2,5-difluorophenyl)thio]ethyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, (-)- (9CI) (CA INDEX NAME)

Rotation (-).

Updated Search

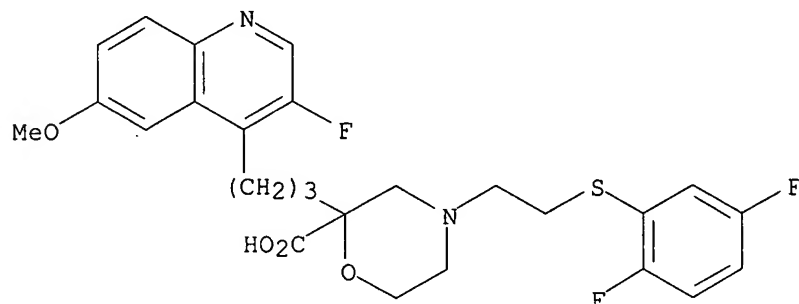
10508761



RN 767355-35-1 HCAPLUS

CN 2-Morpholinecarboxylic acid, 4-[2-[(2,5-difluorophenyl)thio]ethyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, (+)-(9CI) (CA INDEX NAME)

Rotation (+).

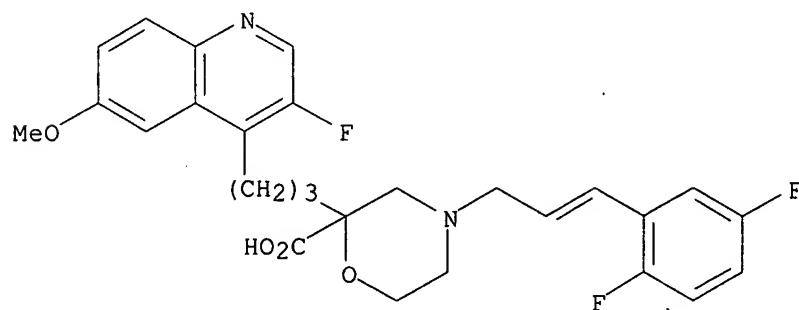


RN 767355-42-0 HCAPLUS

CN 2-Morpholinecarboxylic acid, 4-[3-(2,5-difluorophenyl)-2-propenyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, (-)-(9CI) (CA INDEX NAME)

Rotation (-).

Double bond geometry unknown.



RN 767355-44-2 HCAPLUS

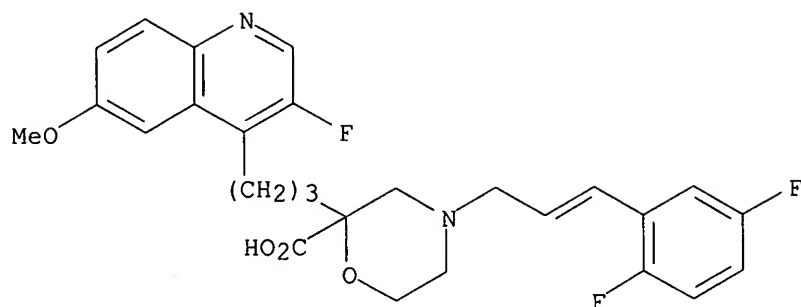
CN 2-Morpholinecarboxylic acid, 4-[3-(2,5-difluorophenyl)-2-propenyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, (+)-(9CI) (CA INDEX NAME)

Rotation (+).

Double bond geometry unknown.

Updated Search

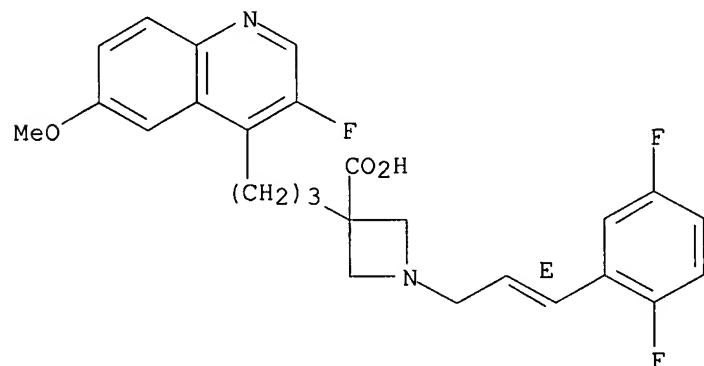
10508761



RN 767355-47-5 HCAPLUS

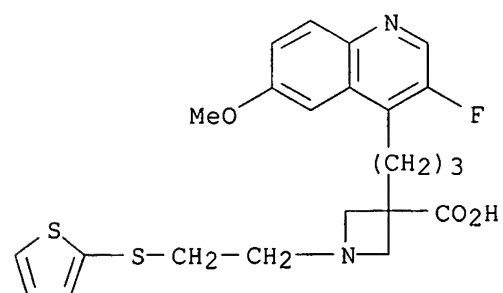
CN 3-Azetidinecarboxylic acid, 1-[(2E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 767355-52-2 HCAPLUS

CN 3-Azetidinecarboxylic acid, 3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-1-[2-(2-thienylthio)ethyl]-, sodium salt (9CI) (CA INDEX NAME)



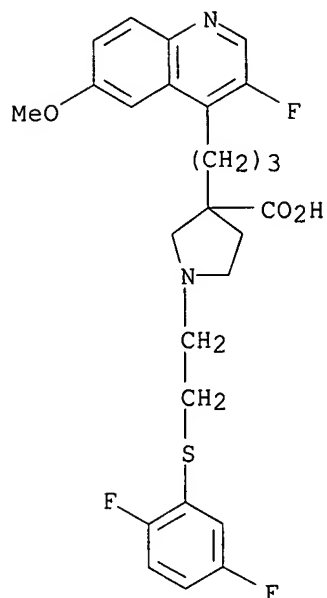
● Na

RN 767355-56-6 HCAPLUS

CN 3-Pyrrolidinecarboxylic acid, 1-[2-[(2,5-difluorophenyl)thio]ethyl]-3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]- (9CI) (CA INDEX NAME)

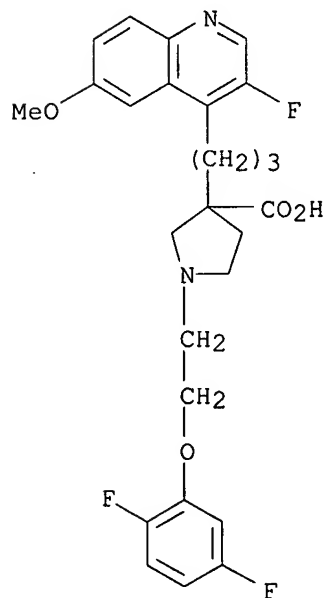
Updated Search

10508761



RN 767355-57-7 HCAPLUS

CN 3-Pyrrolidinecarboxylic acid, 1-[2-(2,5-difluorophenoxy)ethyl]-3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]- (9CI) (CA INDEX NAME)

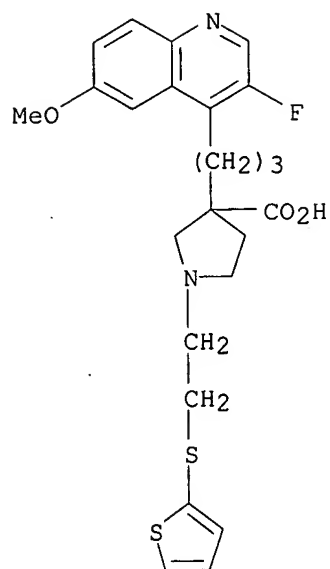


RN 767355-58-8 HCAPLUS

CN 3-Pyrrolidinecarboxylic acid, 3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-1-[2-(2-thienylthio)ethyl]- (9CI) (CA INDEX NAME)

Updated Search

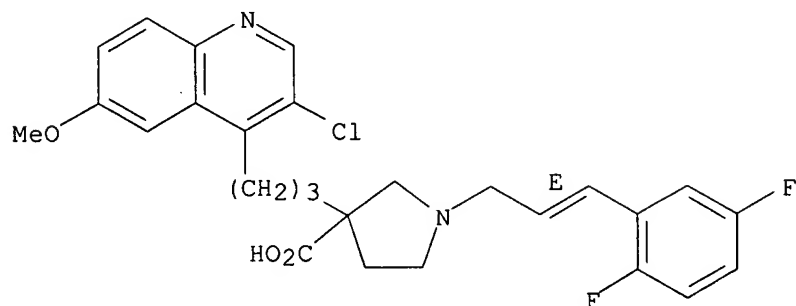
10508761



RN 767355-60-2 HCAPLUS

CN 3-Pyrrolidinecarboxylic acid, 3-[3-(3-chloro-6-methoxy-4-quinolinyl)propyl]-1-[(2E)-3-(2,5-difluorophenyl)-2-propenyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

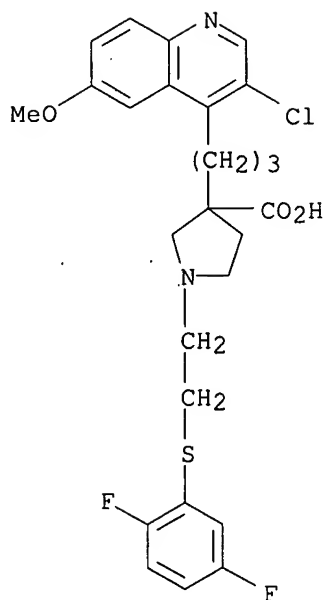


RN 767355-61-3 HCAPLUS

CN 3-Pyrrolidinecarboxylic acid, 3-[3-(3-chloro-6-methoxy-4-quinolinyl)propyl]-1-[2-[(2,5-difluorophenyl)thio]ethyl]- (9CI) (CA INDEX NAME)

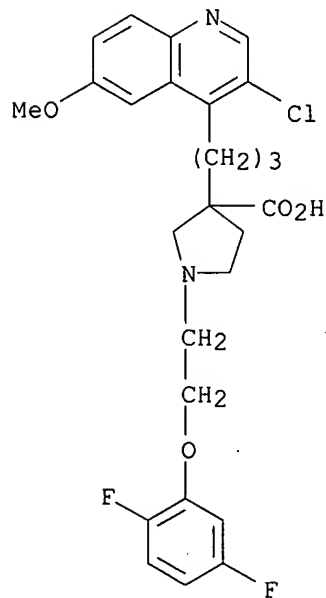
Updated Search

10508761



RN 767355-62-4 HCAPLUS

CN 3-Pyrrolidinecarboxylic acid, 3-[3-(3-chloro-6-methoxy-4-quinolinyl)propyl]-1-[2-(2,5-difluorophenoxy)ethyl]- (9CI) (CA INDEX NAME)

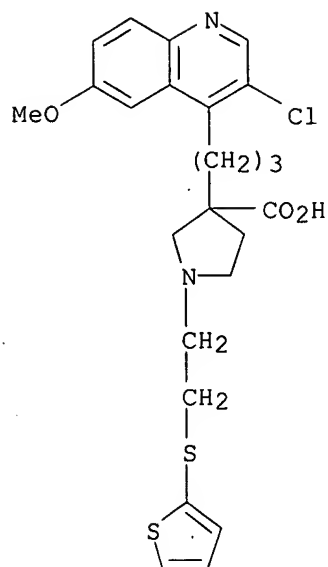


RN 767355-63-5 HCAPLUS

CN 3-Pyrrolidinecarboxylic acid, 3-[3-(3-chloro-6-methoxy-4-quinolinyl)propyl]-1-[2-(2-thienylthio)ethyl]- (9CI) (CA INDEX NAME)

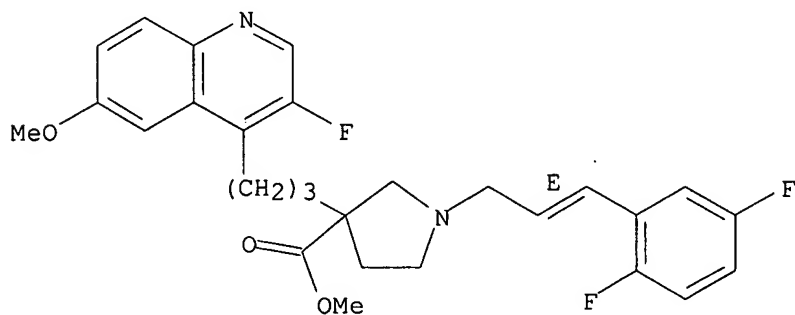
Updated Search

10508761



IT 767355-24-8P, Methyl 1-[(E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylate
767355-38-4P, 2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[2-[(2,5-difluorophenyl)sulfanyl]ethyl]morpholine-2-carboxylic acid methyl ester
RL: PEP (Physical, engineering or chemical process); PYP (Physical process); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); PROC (Process); RACT (Reactant or reagent)
(intermediate; preparation of 4-substituted quinolines as antimicrobials)
RN 767355-24-8 HCAPLUS
CN 3-Pyrrolidinecarboxylic acid, 1-[(2E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester (9CI) (CA INDEX NAME)

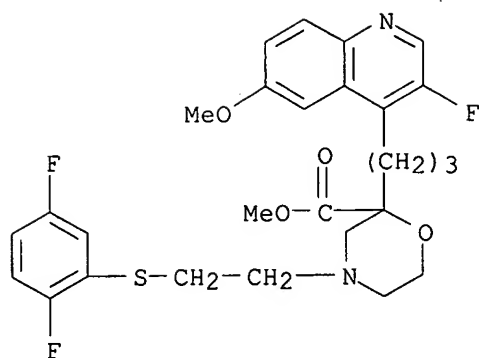
Double bond geometry as shown.



RN 767355-38-4 HCAPLUS
CN 2-Morpholinecarboxylic acid, 4-[2-[(2,5-difluorophenyl)thio]ethyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester (9CI) (CA INDEX NAME)

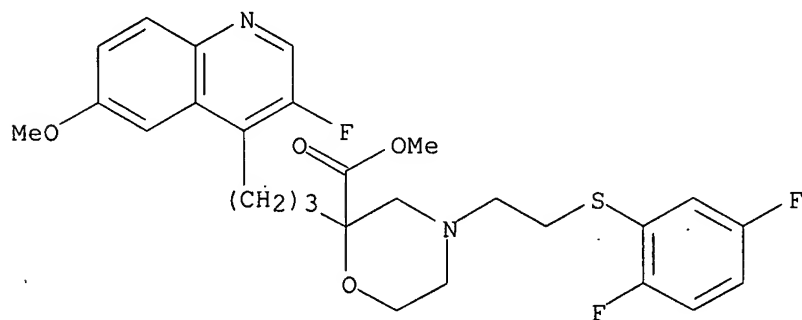
Updated Search

10508761



IT 767355-34-0P, (-)-2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[2-[(2,5-difluorophenyl)sulfanyl]ethyl]morpholine-2-carboxylic acid methyl ester 767355-36-2P, (+)-2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[2-[(2,5-difluorophenyl)sulfanyl]ethyl]morpholine-2-carboxylic acid methyl ester 767355-43-1P, (2R)-2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[3-(2,5-difluorophenyl)-2-propenyl]morpholine-2-carboxylic acid methyl ester 767355-45-3P, (2S)-2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[2-(2,5-difluorophenyl)allyl]morpholine-2-carboxylic acid methyl ester
 RL: PUR (Purification or recovery); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
 (intermediate; preparation of 4-substituted quinolines as antimicrobials)
 RN 767355-34-0 HCAPLUS
 CN 2-Morpholinecarboxylic acid, 4-[2-[(2,5-difluorophenyl)thio]ethyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester, (-)- (9CI) (CA INDEX NAME)

Rotation (-).

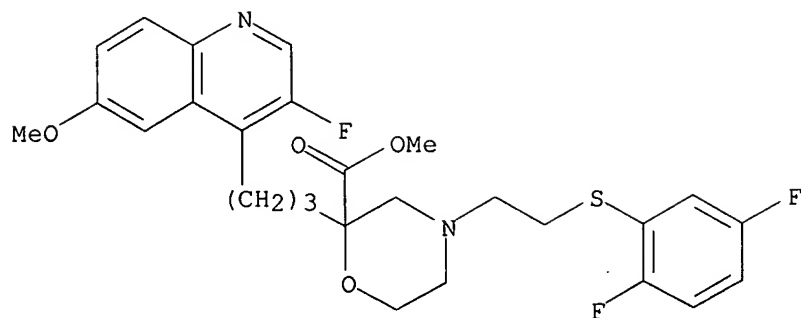


RN 767355-36-2 HCAPLUS
 CN 2-Morpholinecarboxylic acid, 4-[2-[(2,5-difluorophenyl)thio]ethyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester, (+)- (9CI) (CA INDEX NAME)

Rotation (+).

Updated Search

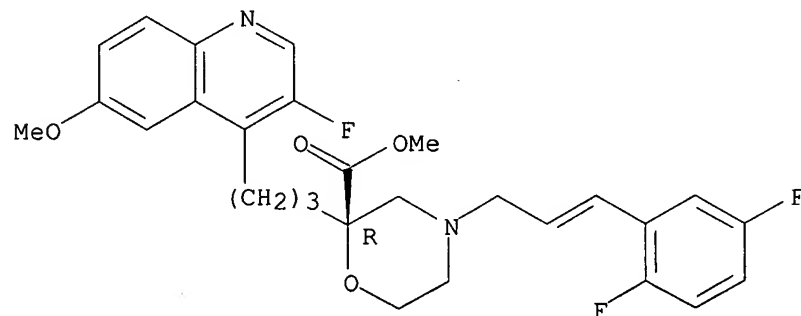
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RN 767355-43-1 HCAPLUS

CN 2-Morpholinecarboxylic acid, 4-[3-(2,5-difluorophenyl)-2-propenyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester, (2R)- (9CI) (CA INDEX NAME)

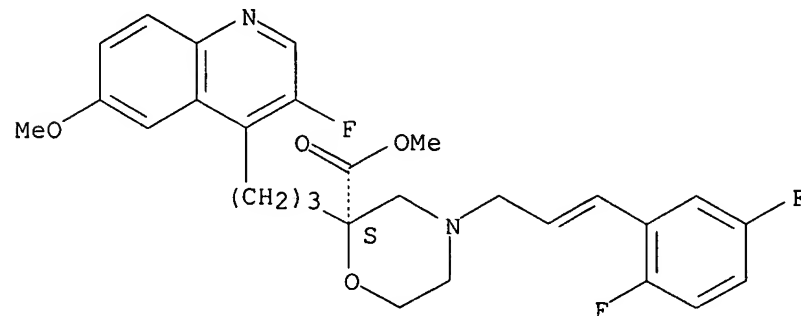
Absolute stereochemistry.
Double bond geometry unknown.



RN 767355-45-3 HCAPLUS

CN 2-Morpholinecarboxylic acid, 4-[3-(2,5-difluorophenyl)-2-propenyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester, (2S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.



IT 767355-25-9P, Methyl 3-[3-(3-fluoro-6-methoxyquinolin-4-

Updated Search

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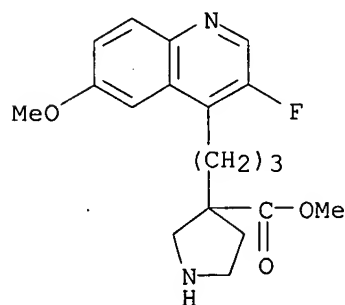
yl)propyl]pyrrolidine-3-carboxylate dihydrochloride 767355-27-1P
, Methyl 1-(tert-butyloxycarbonyl)-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylate 767355-37-3P,
2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]morpholine-2-carboxylic acid
methyl ester 767355-39-5P, 2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-(tert-butyloxycarbonyl)morpholine-2-carboxylic acid methyl
ester 767355-48-6P, Methyl 1-[(E)-3-(2,5-difluorophenyl)-2-
propenyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]azetidine-3-
carboxylate 767355-49-7P, Methyl 3-[3-(3-fluoro-6-
methoxyquinolin-4-yl)propyl]azetidine-3-carboxylate dihydrochloride
767355-50-0P, Methyl 1-(tert-butyloxycarbonyl)-3-[3-(3-fluoro-6-
methoxyquinolin-4-yl)propyl]azetidine-3-carboxylate 767355-53-3P
, Methyl 3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]-1-[2-[(thiophen-2-
yl)sulfanyl]ethyl]azetidine-3-carboxylate 767355-54-4P, Methyl
3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]-1-(2-hydroxyethyl)azetidine-
3-carboxylate

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
(Reactant or reagent)

(intermediate; preparation of 4-substituted quinolines as antimicrobials)

RN 767355-25-9 HCAPLUS

CN 3-Pyrrolidinedicarboxylic acid, 3-[3-(3-fluoro-6-methoxy-4-
quinolinyl)propyl]-, methyl ester, dihydrochloride (9CI) (CA INDEX NAME)

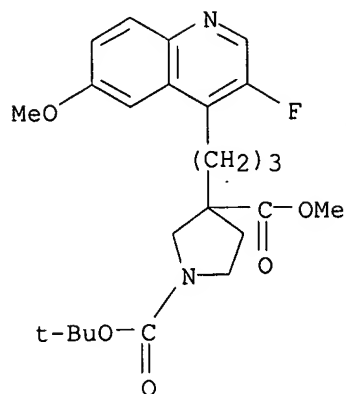


● 2 HCl

RN 767355-27-1 HCAPLUS

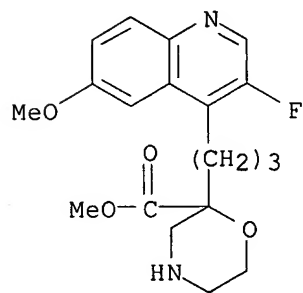
CN 1,3-Pyrrolidinedicarboxylic acid, 3-[3-(3-fluoro-6-methoxy-4-
quinolinyl)propyl]-, 1-(1,1-dimethylethyl) 3-methyl ester (9CI) (CA INDEX
NAME)

10508761



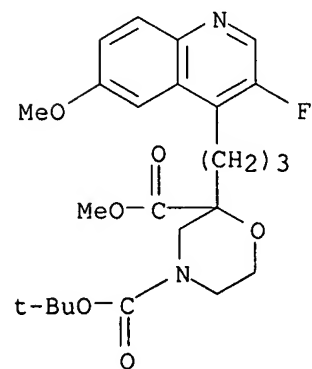
RN 767355-37-3 HCAPLUS

CN 2-Morpholinecarboxylic acid, 2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester (9CI) (CA INDEX NAME)



RN 767355-39-5 HCAPLUS

CN 2,4-Morpholinedicarboxylic acid, 2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, 4-(1,1-dimethylethyl) 2-methyl ester (9CI) (CA INDEX NAME)



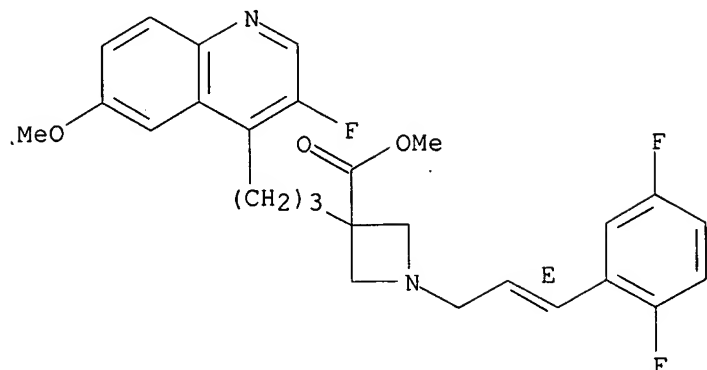
RN 767355-48-6 HCAPLUS

CN 3-Azetidinecarboxylic acid, 1-[(2E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester (9CI) (CA INDEX NAME)

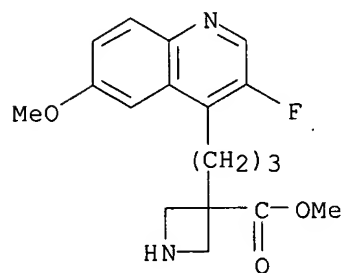
Updated Search

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Double bond geometry as shown.

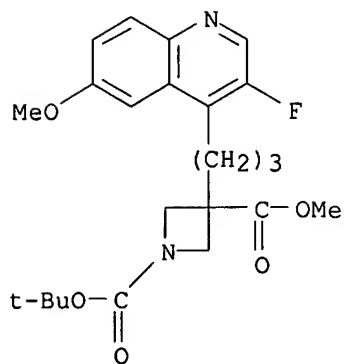


RN 767355-49-7 HCAPLUS
CN 3-Azetidinecarboxylic acid, 3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester, dihydrochloride (9CI) (CA INDEX NAME)



● 2 HCl

RN 767355-50-0 HCAPLUS
CN 1,3-Azetidinedicarboxylic acid, 3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, 1-(1,1-dimethylethyl) 3-methyl ester (9CI) (CA INDEX NAME)

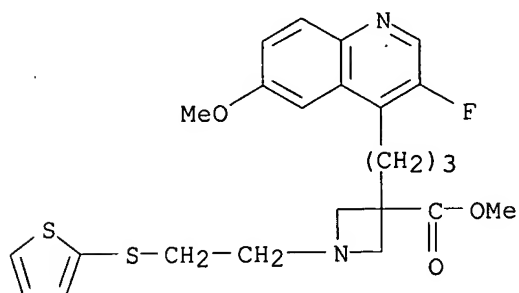


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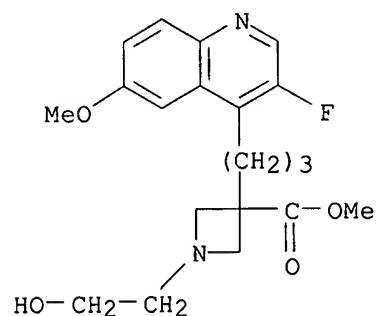
RN 767355-53-3 HCAPLUS

CN 3-Azetidinecarboxylic acid, 3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-1-[2-(2-thienylthio)ethyl]-, methyl ester (9CI) (CA INDEX NAME)



RN 767355-54-4 HCAPLUS

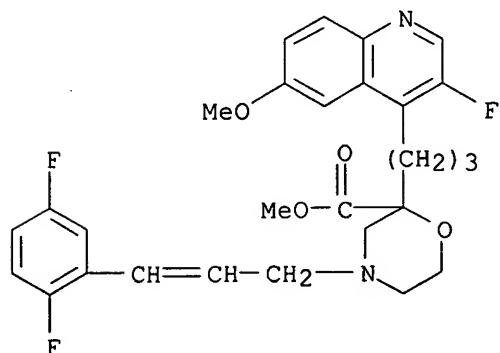
CN 3-Azetidinecarboxylic acid, 3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-1-(2-hydroxyethyl)-, methyl ester (9CI) (CA INDEX NAME)



IT 767355-46-4P, 2-[3-(3-Fluoro-6-methoxyquinolin-4-yl)propyl]-4-[3-(2,5-difluorophenyl)-2-propenyl]morpholine-2-carboxylic acid methyl ester
RL: PEP (Physical, engineering or chemical process); PYP (Physical process); SPN (Synthetic preparation); PREP (Preparation); PROC (Process)
(preparation of 4-substituted quinolines as antimicrobials)

RN 767355-46-4 HCAPLUS

CN 2-Morpholinecarboxylic acid, 4-[3-(2,5-difluorophenyl)-2-propenyl]-2-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester (9CI) (CA INDEX NAME)

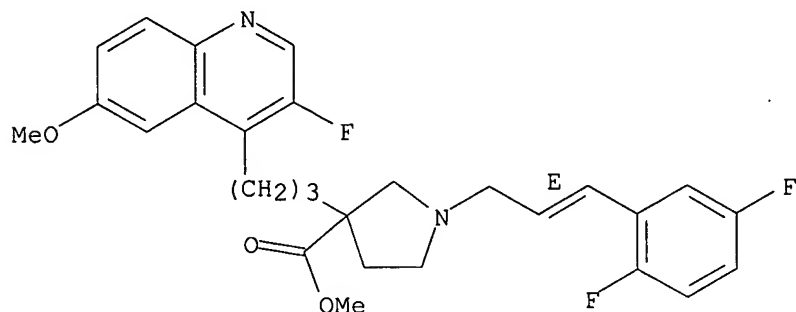


Updated Search

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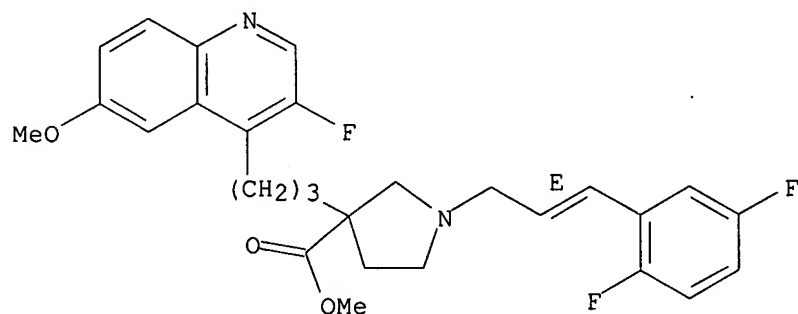
IT 767355-30-6P, (-)-Methyl 1-[(E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylate
767355-32-8P, (+)-Methyl 1-[(E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxyquinolin-4-yl)propyl]pyrrolidine-3-carboxylate
RL: PUR (Purification or recovery); RCT (Reactant); PREP (Preparation);
RACT (Reactant or reagent)
(preparation of 4-substituted quinolines as antimicrobials)
RN 767355-30-6 HCAPLUS
CN 3-Pyrrolidinecarboxylic acid, 1-[(2E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester, (-)- (9CI)
(CA INDEX NAME)

Rotation (-).
Double bond geometry as shown.



RN 767355-32-8 HCAPLUS
CN 3-Pyrrolidinecarboxylic acid, 1-[(2E)-3-(2,5-difluorophenyl)-2-propenyl]-3-[3-(3-fluoro-6-methoxy-4-quinolinyl)propyl]-, methyl ester, (+)- (9CI)
(CA INDEX NAME)

Rotation (+).
Double bond geometry as shown.



REFERENCE COUNT: 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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ENTRY	SESSION
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FULL ESTIMATED COST

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L1	STRUCTURE UPLOADED
L2	0 S L1
L3	37 S L1 FULL

FILE 'HCAPLUS' ENTERED AT 16:53:32 ON 25 JAN 2007

L4	1 S L3
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FILE 'CAOLD' ENTERED AT 16:53:59 ON 25 JAN 2007

=> s l3

L5	0 L3
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=> file reg

	SINCE FILE	TOTAL
	ENTRY	SESSION
COST IN U.S. DOLLARS		
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DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)		
CA SUBSCRIBER PRICE	0.00	-0.78

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L6 STRUCTURE UPLOADED

=> d 16

L6 HAS NO ANSWERS

L6 STR

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Structure attributes must be viewed using STN Express query preparation.

=> s 16

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SAMPLE SCREEN SEARCH COMPLETED - 764414 TO ITERATE

0.3% PROCESSED 2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **INCOMPLETE**

PROJECTED ITERATIONS: 15240514 TO 15336046

PROJECTED ANSWERS: 0 TO 0

L7 0 SEA SSS SAM L6

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L8 STRUCTURE UPLOADED

=> d 18

L8 HAS NO ANSWERS

L8 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

Updated Search

10508761

=> s 18

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1.0% PROCESSED 2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **INCOMPLETE**
PROJECTED ITERATIONS: 4114119 TO 4167441
PROJECTED ANSWERS: 0 TO 0

L9 0 SEA SSS SAM L8

=>

Uploading C:\Documents and Settings\brobinson1\My
Documents\stnweb\Queries\34a541.str

L10 STRUCTURE UPLOADED

=> d 110

L10 HAS NO ANSWERS
L10 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s 110

SAMPLE SEARCH INITIATED 17:08:09 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 207039 TO ITERATE

1.0% PROCESSED 2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **INCOMPLETE**
PROJECTED ITERATIONS: 4114119 TO 4167441
PROJECTED ANSWERS: 0 TO 0

L11 0 SEA SSS SAM L10

=>

Uploading C:\Documents and Settings\brobinson1\My Documents\stnweb\Queries\3434t.str

L12 STRUCTURE UPLOADED

=> d 112

L12 HAS NO ANSWERS
L12 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

Updated Search

10508761

=> s l12

SAMPLE SEARCH INITIATED 17:09:40 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 37636 TO ITERATE

5.3% PROCESSED 2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 741129 TO 764311
PROJECTED ANSWERS: 0 TO 0

L13 0 SEA SSS SAM L12

=> s l12 full

THE ESTIMATED SEARCH COST FOR FILE 'REGISTRY' IS 171.65 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y
FULL SEARCH INITIATED 17:09:45 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 748846 TO ITERATE

100.0% PROCESSED 748846 ITERATIONS
SEARCH TIME: 00.00.06

35 ANSWERS

L14 35 SEA SSS FUL L12

=> file hcaplus

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
183.35	370.73

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
0.00	-0.78

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FILE 'HCAPLUS' ENTERED AT 17:09:54 ON 25 JAN 2007
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FILE COVERS 1907 - 25 Jan 2007 VOL 146 ISS 5
FILE LAST UPDATED: 24 Jan 2007 (20070124/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

Updated Search

10508761

=> s l14
L15 1 L14

=> file reg
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
2.60	373.33

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
0.00	-0.78

CA SUBSCRIBER PRICE

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STRUCTURE FILE UPDATES: 24 JAN 2007 HIGHEST RN 918400-64-3
DICTIONARY FILE UPDATES: 24 JAN 2007 HIGHEST RN 918400-64-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 30, 2006

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conducting SmartSELECT searches.

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predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>
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Documents\stnweb\Queries\55619i.str

L16 STRUCTURE UPLOADED

=> s l16
SAMPLE SEARCH INITIATED 17:12:05 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 207039 TO ITERATE

1.0% PROCESSED 2000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **INCOMPLETE**
PROJECTED ITERATIONS: 4114119 TO 4167441
PROJECTED ANSWERS: 0 TO 0

L17 0 SEA SSS SAM L16

=>

Updated Search

10508761

Uploading C:\Documents and Settings\brobinson1\My Documents\stnweb\Queries\serel.str

L18 STRUCTURE UPLOADED

=> s l18

SAMPLE SEARCH INITIATED 17:12:56 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 206665 TO ITERATE

1.0% PROCESSED 2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **INCOMPLETE**
PROJECTED ITERATIONS: 4106662 TO 4159938
PROJECTED ANSWERS: 0 TO 0

L19 0 SEA SSS SAM L18

=>

Uploading C:\Documents and Settings\brobinson1\My Documents\stnweb\Queries\980u.str

L20 STRUCTURE UPLOADED

=> d l20

L20 HAS NO ANSWERS

L20 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s l20

SAMPLE SEARCH INITIATED 17:14:59 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 20601 TO ITERATE

9.7% PROCESSED 2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 403428 TO 420612
PROJECTED ANSWERS: 0 TO 0

L21 0 SEA SSS SAM L20

=> s l20full

L22 0 L20FULL

=> s l20 full

THE ESTIMATED SEARCH COST FOR FILE 'REGISTRY' IS 171.65 U.S. DOLLARS

DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y

FULL SEARCH INITIATED 17:15:15 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 411764 TO ITERATE

100.0% PROCESSED 411764 ITERATIONS
SEARCH TIME: 00.00.02

35 ANSWERS

Updated Search

10508761

L23 35 SEA SSS FUL L20

=> d his

(FILE 'HOME' ENTERED AT 16:43:46 ON 25 JAN 2007)

FILE 'REGISTRY' ENTERED AT 16:43:54 ON 25 JAN 2007

L1 STRUCTURE UPLOADED

L2 0 S L1

L3 37 S L1 FULL

FILE 'HCAPLUS' ENTERED AT 16:53:32 ON 25 JAN 2007

L4 1 S L3

FILE 'CAOLD' ENTERED AT 16:53:59 ON 25 JAN 2007

L5 0 S L3

FILE 'REGISTRY' ENTERED AT 16:54:06 ON 25 JAN 2007

L6 STRUCTURE UPLOADED

L7 0 S L6

L8 STRUCTURE UPLOADED

L9 0 S L8

L10 STRUCTURE UPLOADED

L11 0 S L10

L12 STRUCTURE UPLOADED

L13 0 S L12

L14 35 S L12 FULL

FILE 'HCAPLUS' ENTERED AT 17:09:54 ON 25 JAN 2007

L15 1 S L14

FILE 'REGISTRY' ENTERED AT 17:10:06 ON 25 JAN 2007

L16 STRUCTURE UPLOADED

L17 0 S L16

L18 STRUCTURE UPLOADED

L19 0 S L18

L20 STRUCTURE UPLOADED

L21 0 S L20

L22 0 S L20FULL

L23 35 S L20 FULL

=> s l23 not l14

L24 0 L23 NOT L14.

=> file reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

180.65

553.98

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

0.00

-0.78

FILE 'REGISTRY' ENTERED AT 17:15:33 ON 25 JAN 2007

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10508761

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 24 JAN 2007 HIGHEST RN 918400-64-3
DICTIONARY FILE UPDATES: 24 JAN 2007 HIGHEST RN 918400-64-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 30, 2006

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>

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L25 STRUCTURE UPLOADED

=> d 125

L25 HAS NO ANSWERS

L25 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s 125

SAMPLE SEARCH INITIATED 17:17:21 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 62502 TO ITERATE

3.2% PROCESSED 2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 1235145 TO 1264935

PROJECTED ANSWERS: 0 TO 0

L26 0 SEA SSS SAM L25

=>

Uploading C:\Documents and Settings\brobinson1\My Documents\stnweb\Queries\2323li.str

L27 STRUCTURE UPLOADED

=> d 127

L27 HAS NO ANSWERS

L27 STR

Updated Search

10508761

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *
Structure attributes must be viewed using STN Express query preparation.

=> s 127

SAMPLE SEARCH INITIATED 17:18:15 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 19665 TO ITERATE

10.2% PROCESSED 2000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 384905 TO 401695
PROJECTED ANSWERS: 0 TO 0

L28 0 SEA SSS SAM L27

=> s 127 full

THE ESTIMATED SEARCH COST FOR FILE 'REGISTRY' IS 171.65 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y
FULL SEARCH INITIATED 17:18:20 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 391797 TO ITERATE

100.0% PROCESSED 391797 ITERATIONS 35 ANSWERS
SEARCH TIME: 00.00.02

L29 35 SEA SSS FUL L27

=>

Uploading C:\Documents and Settings\brobinson1\My Documents\stnweb\Queries\1121j.str

L30 STRUCTURE UPLOADED

=> s 130

SAMPLE SEARCH INITIATED 17:19:51 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 45759 TO ITERATE

4.4% PROCESSED 2000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 902411 TO 927949
PROJECTED ANSWERS: 0 TO 0

L31 0 SEA SSS SAM L30

=> s 130 full

THE ESTIMATED SEARCH COST FOR FILE 'REGISTRY' IS 171.65 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y
FULL SEARCH INITIATED 17:19:56 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 911515 TO ITERATE

100.0% PROCESSED 911515 ITERATIONS 35 ANSWERS

Updated Search

10508761

SEARCH TIME: 00.00.03

L32 35 SEA SSS FUL L30

=>

Uploading C:\Documents and Settings\brobinson1\My
Documents\stnweb\Queries\2323km.str

L33 STRUCTURE UPLOADED

=>

Uploading C:\Documents and Settings\brobinson1\My
Documents\stnweb\Queries\121a2k.str

L34 STRUCTURE UPLOADED

=> s 134

SAMPLE SEARCH INITIATED 17:23:49 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 2858 TO ITERATE

70.0% PROCESSED 2000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 53954 TO 60366
PROJECTED ANSWERS: 0 TO 0

L35 0 SEA SSS SAM L34

=> s 134 full

THE ESTIMATED SEARCH COST FOR FILE 'REGISTRY' IS 171.65 U.S. DOLLARS
DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y
FULL SEARCH INITIATED 17:23:56 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 56986 TO ITERATE

100.0% PROCESSED 56986 ITERATIONS (2 INCOMPLETE)
SEARCH TIME: 00.00.05

6 ANSWERS

L36 6 SEA SSS FUL L34

=> file hcaplus

COST IN U.S. DOLLARS

SINCE FILE
ENTRY

TOTAL
SESSION

FULL ESTIMATED COST

521.25

1075.23

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE
ENTRY

TOTAL
SESSION

CA SUBSCRIBER PRICE

0.00

-0.78

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FILE COVERS 1907 - 25 Jan 2007 VOL 146 ISS 5
FILE LAST UPDATED: 24 Jan 2007 (20070124/ED)

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=> s 136
L37

5 L36

=> d 137, ibib abs hitstr, 1-5

L37 ANSWER 1 OF 5 HCAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 2005:1289687 HCAPLUS

DOCUMENT NUMBER: 144:51568

TITLE: Preparation of substituted 2-quinolyl-oxazoles and their heterocyclic analogs useful as pde4 inhibitors
INVENTOR(S): Kuang, Rongze; Blythin, David; Shih, Neng-Yang; Shue, Ho-Jane; Chen, Xiao; Cao, Jianhua; Gu, Danlin; Huang, Ying; Schwerdt, John H.; Ting, Pauline C.; Wong, Shing-Chun; Xiao, Li

PATENT ASSIGNEE(S): Schering Corporation, USA

SOURCE: PCT Int. Appl., 233 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

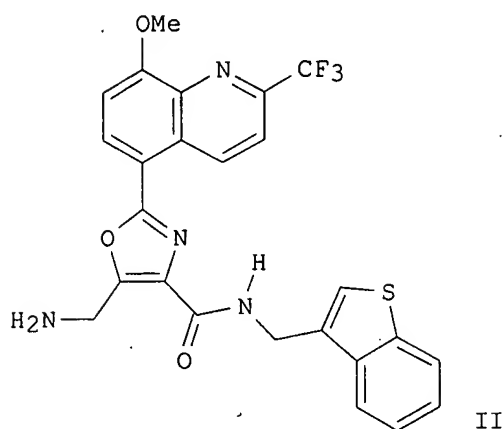
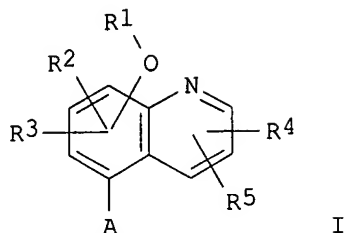
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
-----	----	-----	-----	-----
WO 2005116009	A1	20051208	WO 2005-US17134	20050516
WO 2005116009	B1	20060126		
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KM, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NG, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SM, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
RW:	BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
AU 2005247906	A1	20051208	AU 2005-247906	20050516
CA 2565599	A1	20051208	CA 2005-2565599	20050516
US 2006106062	A1	20060518	US 2005-130359	20050516
PRIORITY APPLN. INFO.:			US 2004-572266P	P 20040518
			WO 2005-US17134	W 20050516

Updated Search

10508761

OTHER SOURCE(S):
GI

MARPAT 144:51568



AB Title compds. I [R1 = H, alkyl, cycloalkyl; R2, R3 and R5 independently = H or halo; R4 = H, halo, alkyl, etc.; A = substituted oxazolyl, imidazole, thiazole or pyrrole], and their pharmaceutically acceptable salts, are prepared and disclosed as pde4 inhibitors. Thus, e.g., II was prepared in a multistep synthesis from 2-trifluoromethyl-8-methoxyquinolin-5-yl carboxylic acid. In PDE4 assays, selected compds. possessed IC50 values ranging from 0.01-1.8 nM. Also claimed are pharmaceutical compns., the use of the compds. as PDE4 inhibitors, and combinations with other actives.

IT 871011-97-1P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

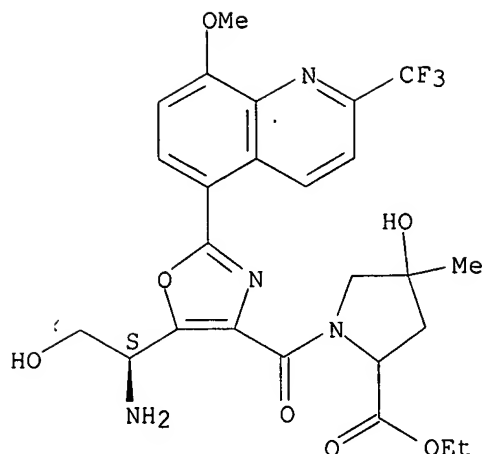
(preparation of substituted quinolyloxazoles and their heterocyclic analogs useful as PDE4 inhibitors)

RN 871011-97-1 HCAPLUS

CN Proline, 1-[[5-[(1S)-1-amino-2-hydroxyethyl]-2-[8-methoxy-2-(trifluoromethyl)-5-quinolinyl]-4-oxazolyl]carbonyl]-4-hydroxy-4-methyl-, ethyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Updated Search



● HCl

REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L37 ANSWER 2 OF 5 HCAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 2004:310829 HCAPLUS

DOCUMENT NUMBER: 140:303552

TITLE: Preparation of β -amino acid derivatives as inhibitors of matrix metalloproteases and TNF- α

INVENTOR(S): Duan, Jingwu; King, Bryan W.; Decicco, Carl; Maduskuie, Thomas P.; Voss, Mathew E.

PATENT ASSIGNEE(S): USA

SOURCE: U.S. Pat. Appl. Publ., 150 pp.

CODEN: USXXCO

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2004072802	A1	20040415	US 2002-267207	20021009
PRIORITY APPLN. INFO.:			US 2002-267207	20021009

OTHER SOURCE(S): MARPAT 140:303552

AB Novel β -amino acid derivs. A-CR3R4aCR2R4NR1CO-X-Z-Ua-Xa-Ya-Za [A = CO₂H, SH, CH₂SH, S(O)Ra:NH (Ra = H, alkyl), P(O)(OH)₂, etc.; X, Xa is absent or alkylene, alkenylene or alkynylene; Z is absent or substituted C3-13 carbocycle or 5-14 membered heterocycle; Ua is absent or O, NRa1 [Ra1 = H, (un)substituted alkyl, alkenyl or alkynyl; Ra and Ra1 may form a ring], CO, CO₂, O₂C, CONRa1, S(O)p (p = 0-2), etc.; Ya is absent or O, NRa1, S(O)p or CO; Za is H, substituted C3-13 carbocycle or 5-14 membered heterocycle; R1 is H, alkyl, Ph, benzyl; R2 is Q (Q is H, substituted carbocycle or heterocycle), alkylene-Q, (CRaRa1)r1O(CRaRa1)r-Q (r, r1 = 0-4), (CRaRa1)r1NRa(CRaRa1)r-Q, etc.; R3 = Q1 (Q1 is any group given for Q), alkylene-Q1, (CRaRa1)r1O(CRaRa1)r-Q1, (CRaRa1)r1NRa(CRaRa1)r-Q1, etc.; R4, R4a = H, substituted alkyl, alkenyl or alkynyl; alternatively R1 and R2, R1 and R3, R3 and R4a may form rings (with provisos)] or a

10508761

stereoisomer or pharmaceutically acceptable salt were prepared as metalloprotease and TNF- α inhibitors. Thus, N-hydroxy-1-[[4-[(2-methyl-4-quinolinyl)methoxy]phenyl]acetyl]-3-azetidinecarboxamide was prepared by a multistep procedure involving reactions of Me 4-hydroxyphenylacetate, 2-methyl-4-quinolinylmethanol, and 3-azetidinecarboxylic acid Me ester.

IT 362700-34-3P 362700-35-4P

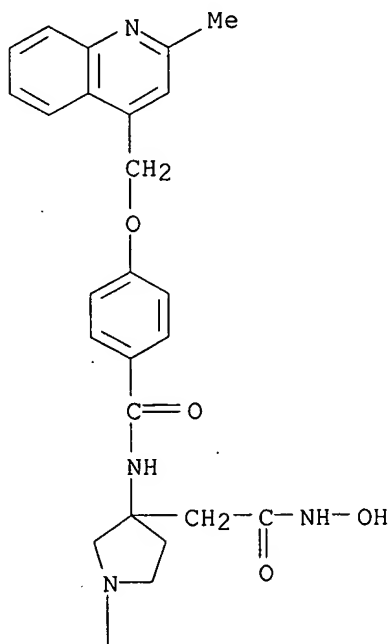
RL: BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of β -amino acid derivs. as inhibitors of matrix metalloproteases and TNF- α)

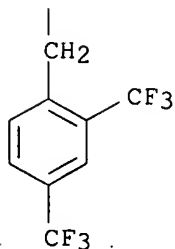
RN 362700-34-3 HCAPLUS

CN 3-Pyrrolidineacetamide, 1-[[2,4-bis(trifluoromethyl)phenyl]methyl]-N-hydroxy-3-[[4-[(2-methyl-4-quinolinyl)methoxy]benzoyl]amino]- (9CI) (CA INDEX NAME)

PAGE 1-A



PAGE 2-A



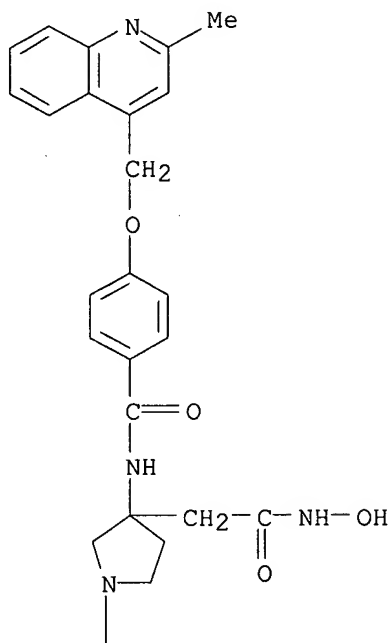
10508761

RN 362700-35-4 HCAPLUS
CN 3-Pyrrolidineacetamide, 1-[[[2,4-bis(trifluoromethyl)phenyl]methyl]-N-hydroxy-3-[[4-[(2-methyl-4-quinolinyl)methoxy]benzoyl]amino]-, bis(trifluoroacetate) (salt) (9CI) (CA INDEX NAME)

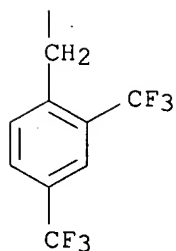
CM 1

CRN 362700-34-3
CMF C33 H30 F6 N4 O4

PAGE 1-A



PAGE 2-A

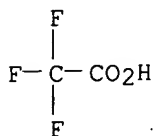


CM 2

CRN 76-05-1
CMF C2 H F3 O2

Updated Search

10508761



L37 ANSWER 3 OF 5 HCAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 2003:616857 HCAPLUS

DOCUMENT NUMBER: 139:286905

TITLE: Efficient incorporation of positively charged 2', 3'-dideoxynucleoside-5'-triphosphates by DNA polymerases and their application in direct-load' DNA sequencing

AUTHOR(S): Finn, Patrick J.; Bull, Matthew G.; Xiao, Haiguang; Phillips, Paula D.; Nelson, John R.; Grossmann, Greg; Nampalli, Satyam; McArdle, Bernard F.; Mamone, J. Anthony; Flick, Parke K.; Fuller, Carl W.; Kumar, Shiv
CORPORATE SOURCE: Amersham Biosciences, Piscataway, NJ, 08855-1327, USA
SOURCE: Nucleic Acids Research (2003), 31(16), 4769-4778

CODEN: NARHAD; ISSN: 0305-1048

PUBLISHER: Oxford University Press

DOCUMENT TYPE: Journal

LANGUAGE: English

AB A series of charge-modified, dye-labeled 2', 3'-dideoxynucleoside-5'-triphosphates have been synthesized and evaluated as reagents for dye-terminator DNA sequencing. Unlike the commonly used dye-labeled terminators, these terminators possess a net pos. charge and migrate in the opposite direction to dye-labeled Sanger fragments during electrophoresis. Post-sequencing reaction purification is not required to remove unreacted nucleotide or associated breakdown products prior to electrophoresis. Thus, DNA sequencing reaction mixts. can be loaded directly onto a separating medium such as a sequencing gel. The charge-modified nucleotides have also been shown to be more efficiently incorporated by a number of DNA polymerases than regular dye-labeled dideoxynucleotide terminators or indeed normal dideoxynucleoside-5'-triphosphates.

IT 608520-72-5P

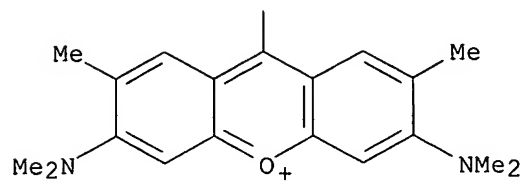
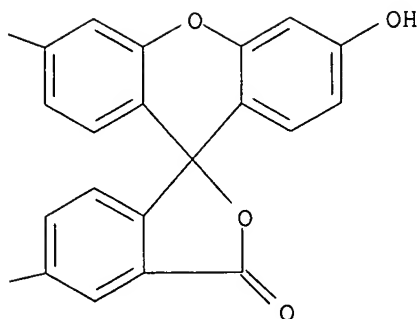
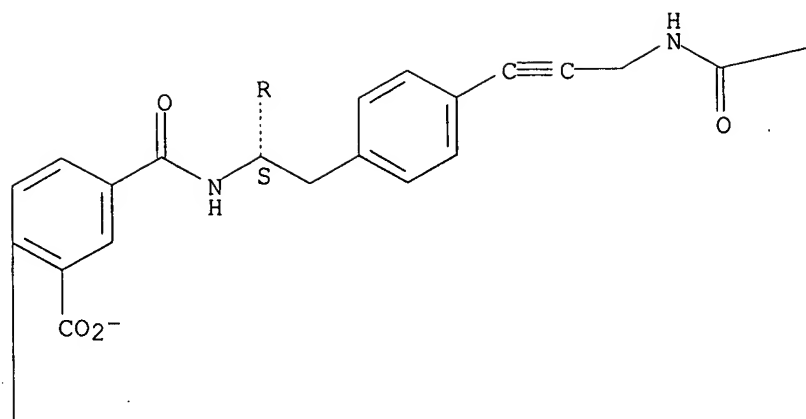
RL: ARG (Analytical reagent use); SPN (Synthetic preparation); ANST (Analytical study); PREP (Preparation); USES (Uses)
(efficient incorporation of pos. charged 2', 3'-dideoxynucleoside-5'-triphosphates by DNA polymerases and their application in direct-load' DNA sequencing)

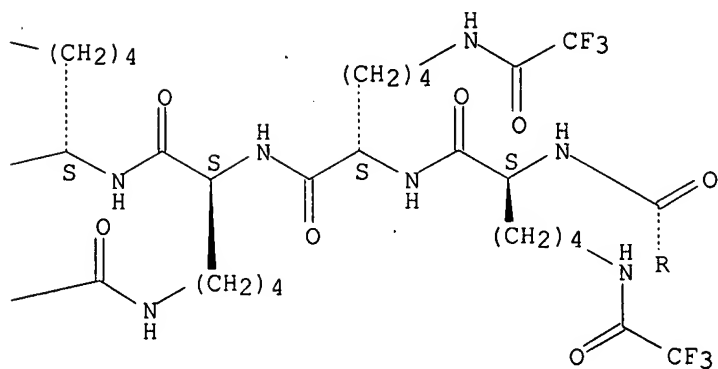
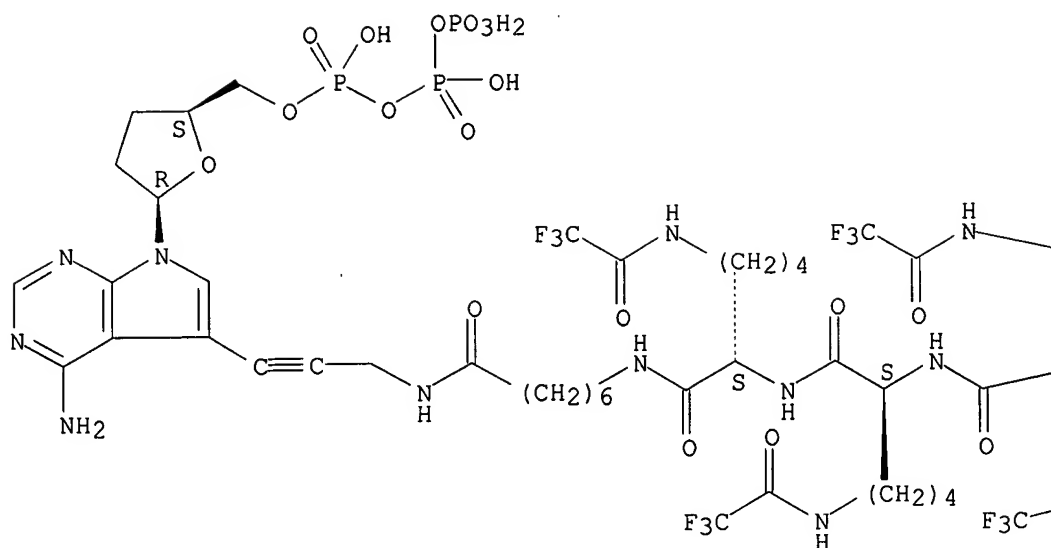
RN 608520-72-5 HCAPLUS

CN L-Lysinamide, N-[4-[3,6-bis(dimethylamino)-2,7-dimethylxanthylum-9-yl]-3-carboxybenzoyl]-4-[3-[[[3',6'-dihydroxy-3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthen]-5-yl]carbonyl]amino]-1-propynyl]-L-phenylalanyl-N6-(trifluoroacetyl)-L-lysyl-N6-(trifluoroacetyl)-L-lysyl-N6-(trifluoroacetyl)-L-lysyl-N6-(trifluoroacetyl)-L-lysyl-N6-(trifluoroacetyl)-L-lysyl-N-[7-[[3-[4-amino-7-[(2R,5S)-tetrahydro-5-(3,5,7,7-tetrahydroxy-3,5,7-trioxido-2,4,6-trioxa-3,5,7-triphosphahept-1-yl)-2-furanyl]-7H-pyrrolo[2,3-d]pyrimidin-5-yl]-2-propynyl]amino]-7-oxoheptyl]-N6-(trifluoroacetyl)-, inner salt (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Updated Search





REFERENCE COUNT: 16 THERE ARE 16 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L37 ANSWER 4 OF 5 HCAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 2001:713343 HCAPLUS

DOCUMENT NUMBER: 135:272894

TITLE: Preparation of β -amino acid derivatives as inhibitors of matrix metalloproteases and TNF- α

INVENTOR(S): Duan, Jingwu; King, Bryan W.; Decicco, Carl; Maduskuie, Thomas P., Jr.; Voss, Matthew E.

10508761

PATENT ASSIGNEE(S): Dupont Pharmaceuticals Company, USA
 SOURCE: PCT Int. Appl., 483 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001070734	A2	20010927	WO 2001-US8336	20010315
WO 2001070734	A3	20020314		
W: AT, AU, BR, CA, CH, CN, CZ, DE, DK, EE, ES, FI, GB, HU, IL, IN, JP, KR, LT, LU, LV, NZ, PL, PT, RO, SE, SG, SI, SK, UA, VN, ZA, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR				
CA 2400168	A1	20010927	CA 2001-2400168	20010315
AU 200150850	A	20011003	AU 2001-50850	20010315
EP 1263756	A2	20021211	EP 2001-924171	20010315
EP 1263756	B1	20040225		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, CY, TR				
BR 2001009469	A	20030429	BR 2001-9469	20010315
JP 2003528097	T	20030924	JP 2001-568935	20010315
AT 260272	T	20040315	AT 2001-924171	20010315
NZ 521245	A	20040430	NZ 2001-521245	20010315
ES 2215893	T3	20041016	ES 2001-1924171	20010315
US 2002013341	A1	20020131	US 2001-811116	20010316
US 6495565	B2	20021217		
IN 2002MN01075	A	20050304	IN 2002-MN1075	20020808
HK 1049334	A1	20040716	HK 2003-101437	20030226
PRIORITY APPLN. INFO.:				
			US 2000-190183P	P 20000317
			US 2000-235467P	P 20000926
			US 2000-252062P	P 20001120
			WO 2001-US8336	W 20010315

OTHER SOURCE(S): MARPAT 135:272894

AB Novel β -amino acid derivs. A-CR3R4aCR2R4NR1CO-X-Z-Ua-Xa-Ya-Za [A = CO2H, SH, CH2SH, S(O)Ra:NH (Ra = H, alkyl), P(O)(OH)2, etc.; X, Xa is absent or alkylene, alkenylene or alkynylene; Z is absent or substituted C3-13 carbocycle or 5-14 membered heterocycle; Ua is absent or O, NRa1 [Ra1 = H, (un)substituted alkyl, alkenyl or alkynyl; Ra and Ra1 may form a ring], CO, CO2, O2C, CONRa1, S(O)p (p = 0-2), etc.; Ya is absent or O, NRa1, S(O)p or CO; Za is H, substituted C3-13 carbocycle or 5-14 membered heterocycle; R1 is H, alkyl, Ph, benzyl; R2 is Q (Q is H, substituted carbocycle or heterocycle), alkylene-Q, (CRaRa1)r1O(CRaRa1)r-Q (r, r1 = 0-4), (CRaRa1)r1NRa(CRaRa1)r-Q, etc.; R3 = Q1 (Q1 is any group given for Q), alkylene-Q1, (CRaRa1)r1O(CRaRa1)r-Q1, (CRaRa1)r1NRa(CRaRa1)r-Q1, etc.; R4, R4a = H, substituted alkyl, alkenyl or alkynyl; alternatively R1 and R2, R1 and R3, R3 and R4a may form rings (with provisos)] or a stereoisomer or pharmaceutically acceptable salt were prepared as metalloprotease and TNF- α inhibitors. Thus, N-hydroxy-1-[[4-[(2-methyl-4-quinolinyl)methoxy]phenyl]acetyl]-3-azetidinedicarboxamide was prepared by a multistep procedure involving reactions of Me 4-hydroxyphenylacetate, 2-methyl-4-quinolinylmethanol, and 3-azetidinedicarboxylic acid Me ester.

IT 362700-34-3P 362700-35-4P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);

Updated Search

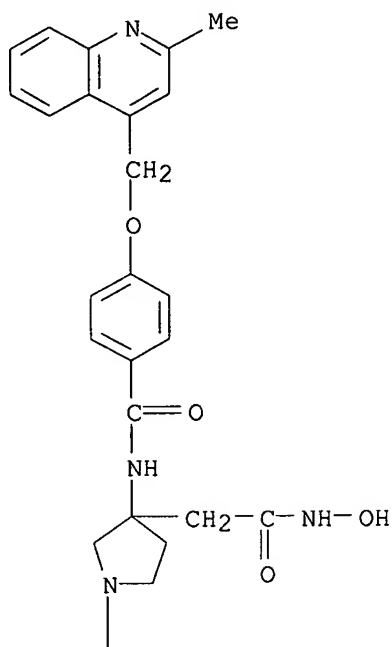
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BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of β -amino acid derivs. as inhibitors of matrix metalloproteases and TNF- α)

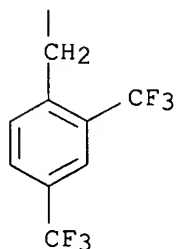
RN 362700-34-3 HCAPLUS

CN 3-Pyrrolidineacetamide, 1-[[2,4-bis(trifluoromethyl)phenyl]methyl]-N-hydroxy-3-[[4-[(2-methyl-4-quinolinyl)methoxy]benzoyl]amino]- (9CI) (CA INDEX NAME)

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RN 362700-35-4 HCAPLUS

CN 3-Pyrrolidineacetamide, 1-[[2,4-bis(trifluoromethyl)phenyl]methyl]-N-hydroxy-3-[[4-[(2-methyl-4-quinolinyl)methoxy]benzoyl]amino]-, bis(trifluoroacetate) (salt) (9CI) (CA INDEX NAME)

CM 1

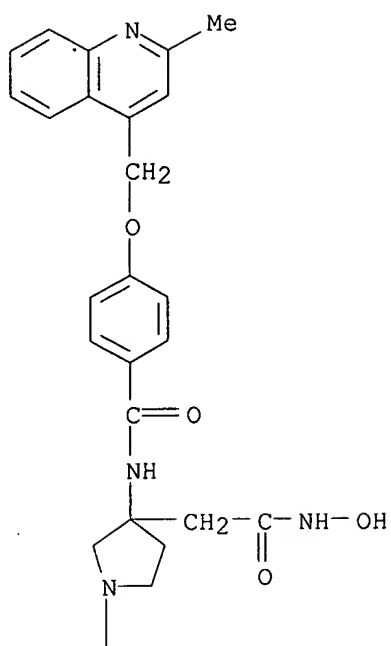
CRN 362700-34-3

Updated Search

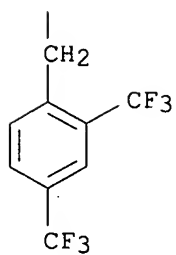
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CMF C33 H30 F6 N4 O4

PAGE 1-A

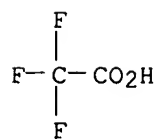


PAGE 2-A



CM 2

CRN 76-05-1
CMF C2 H F3 O2



Updated Search

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L37 ANSWER 5 OF 5 HCAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 2001:380438 HCAPLUS
DOCUMENT NUMBER: 135:24657
TITLE: Selective cellular targeting: multifunctional delivery vehicles
INVENTOR(S): Glazier, Arnold
PATENT ASSIGNEE(S): Drug Innovation & Design, Inc., USA
SOURCE: PCT Int. Appl., 981 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001036003	A2	20010525	WO 2000-US31262	20001114
W:			AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM	
RW:			GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG	
CA 2391534	A1	20010525	CA 2000-2391534	20001114
AU 2001016075	A5	20010530	AU 2001-16075	20001114
EP 1255567	A1	20021113	EP 2000-978631	20001114
R:			AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR	
US 2003138432	A1	20030724	US 2000-738625	20001215
PRIORITY APPLN. INFO.:			US 1999-165485P	P 19991115
			US 2000-239478P	P 20001011
			US 2000-241937P	P 20001020
			WO 2000-US31262	W 20001114
			US 2000-712465	B1 20001115

AB The present invention relates to the compns., methods, and applications of a novel approach to selective cellular targeting. The purpose of this invention is to enable the selective delivery and/or selective activation of effector mols. to target cells for diagnostic or therapeutic purposes. The present invention relates to multi-functional prodrugs or targeting vehicles wherein each functionality is capable of enhancing targeting selectivity, affinity, intracellular transport, activation or detoxification. The present invention also relates to ultralow dose, multiple target, multiple drug chemotherapy and targeted immunotherapy for cancer treatment.

IT 341553-47-7P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); PNU (Preparation, unclassified); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(multifunctional delivery vehicles for selective cellular targeting of drugs)

RN 341553-47-7 HCAPLUS

CN 10,13,16,26,29,32,35,45,48,51-Decaoxa-2,7,19,23,42,54-hexaaza-58-phosphadohexacontane-3,60,62-tricarboxylic acid, 1-[4-[[[(2-amino-1,4-dihydro-4-oxo-6-pteridinyl)methyl][[3-[[[3-carboxy-1-oxopropoxy)methoxy]carbonyl]-4-[(2-oxido-1,3,2-dioxaphosphorinan-2-

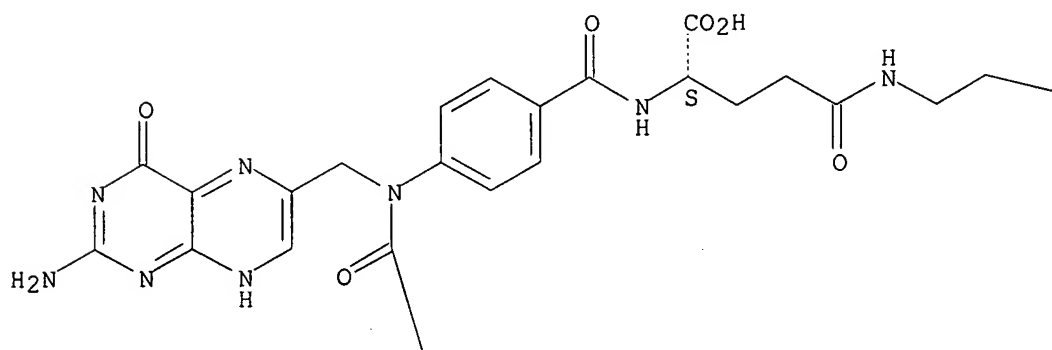
Updated Search

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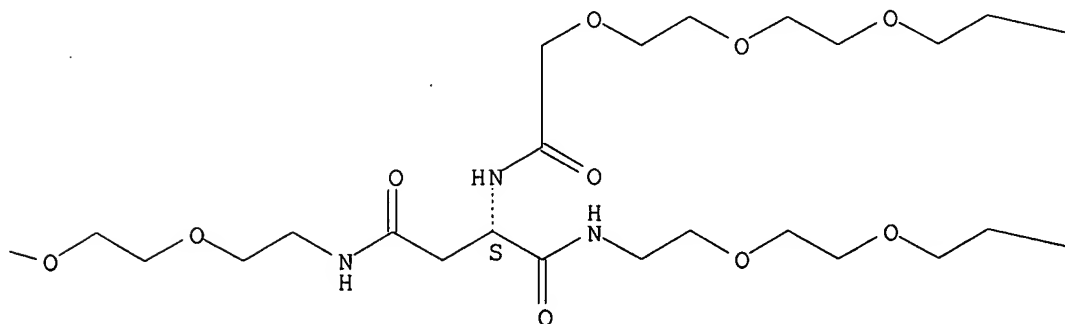
yl)oxy]phenyl]methoxy]carbonyl]amino]phenyl]-22-[17-[6-[[[[[1-[[5-(5-carboxy-3-methyl-2-pentenyl)-1,3-dihydro-6-methoxy-7-methyl-3-oxo-4-isobenzofuranyl]oxy]-2,2,2-trifluoroethyl]amino]carbonyl]oxy]methyl]-5,8-dioxo-1-naphthalenyl]-1,14-dioxo-5,8,11-trioxa-2,15-diazaheptadec-1-yl]-39-(1,14-dioxo-5,8,11-trioxa-2,15-diazadocos-1-yl)-58-hydroxy-1,6,20,24,37,41,55-heptaoso-, 58-oxide, (3S,22S,39S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.

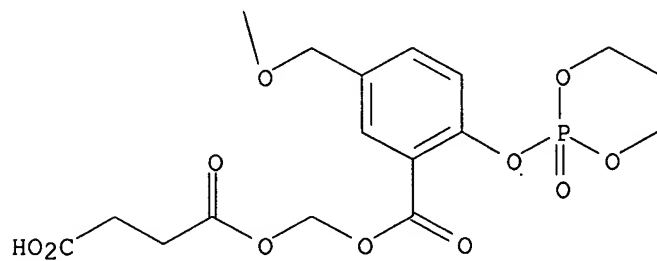
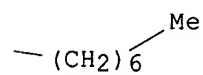
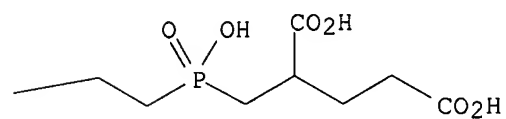
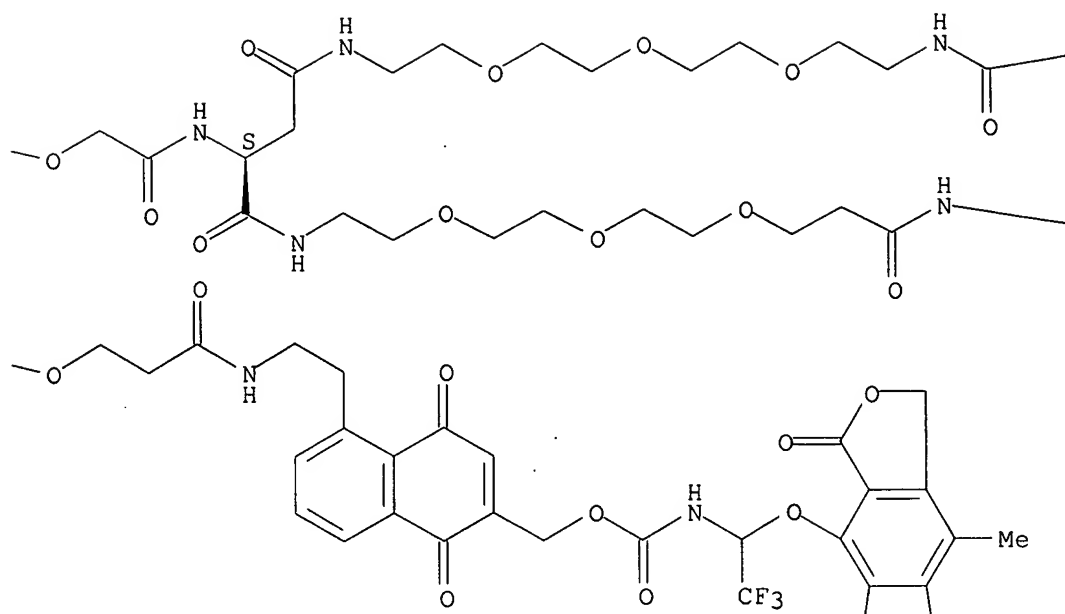
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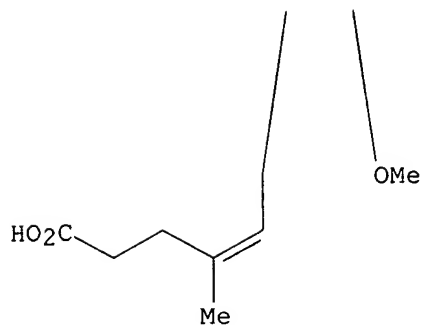


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